

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

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CERTIFICATE

This is to certify that the Lab work entitled “ Bio Inspired Systems (23CS5BSBIS)” carried out by **Sindhuja Narasimhan (1BM22CS279)**, 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/sindhuja279/BISlab.git>

Program 1

Genetic Algorithm for Optimization Problems

Algorithm:

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Lab 2.

Algorithm for genetic algorithm

- Step 1. Initialize parameters: set the population size, mutation rate, crossover rate & no. of generations.
- Step 2. Generate Initial population: create a random population of potential solutions within the given bounds.
- Step 3. Evaluate Fitness: calculate the fitness of each individual in the population by evaluating the objective function.
- Step 4. Select Parents: select the fittest individuals from the population to reproduce based on their fitness.
- Step 5. Crossover: perform crossover between the selected parents to create new offspring, with a probability equal to the crossover rate.
- Step 6. Mutation: apply mutation to the offspring, with a probability equal to mutation rate, to introduce new traits.
- Step 7. Replace Least Fit: replace the least fit individuals in the population with the new offspring.
- Step 8. Repeat: repeat steps 3-7 for a fixed number of generations or until convergence criteria are met.
- Step 9. Output Best Soln: return the best soln found during the generations, which is the individual with the

→ The basic idea behind a genetic algorithm is to simulate the process of evolution by iteratively selecting, breeding and mutating a population of candidate solution.

→ The genetic algorithm is used for

- Optimizing problem such as scheduling, resource allocation and portfolio optimization
- machine learning to optimize parameters of models such as neural networks and decision trees
- engineering design: to optimize complex systems such as electronic circuits etc.
- computer network: to optimize network routing, scheduling and resource allocation.

→ Optimizing Techniques:

- selection: selecting the fittest individuals from the population to reproduce.
- crossover: combining the genetic information of two parents to create a new offspring
- mutation: Randomly changing the genetic information of an individual to introduce new traits.
- elitism: preserving the best solution from the previous generation to ensure that the best solution are not lost.
- Simulated binary crossover: A crossover technique that simulates the process of binary crossover to create new offspring

10/24

Code:

```
import random
import numpy as np

def objective_function(x):
    # Define the objective function to be minimized
    return x**2 + 2*x + 1

def generate_initial_population(population_size, bounds):
    # Generate an initial population of potential solutions
    population = []
    for _ in range(population_size):
        x = random.uniform(bounds[0], bounds[1])
        population.append(x)
    return population

def evaluate_fitness(population):
    # Evaluate the fitness of each individual in the population
    fitness = []
    for x in population:
        fitness.append(objective_function(x))
    return fitness

def selection(population, fitness, num_parents):
    # Select individuals based on their fitness to reproduce
    parents = []
    for _ in range(num_parents):
        max_fitness_idx = np.argmax(fitness)
        parents.append(population[max_fitness_idx])
        fitness[max_fitness_idx] = -float('inf') # Set fitness to -inf to avoid re-selection
    return parents

def crossover(parents, crossover_rate):
    # Perform crossover between selected individuals to produce offspring
    offspring = []
    for _ in range(len(parents) // 2):
        parent1, parent2 = random.sample(parents, 2)
        if random.random() < crossover_rate:
            x1, x2 = parent1, parent2
            x_offspring = (x1 + x2) / 2
            offspring.append(x_offspring)
    return offspring

def mutation(offspring, mutation_rate, bounds):
    # Apply mutation to the offspring to maintain genetic diversity
    for i in range(len(offspring)):
```

```

    if random.random() < mutation_rate:
        offspring[i] += random.uniform(-0.1, 0.1) * (bounds[1] - bounds[0])
        offspring[i] = max(bounds[0], min(offspring[i], bounds[1])) # Ensure bounds
    return offspring

def genetic_algorithm(population_size, mutation_rate, crossover_rate, num_generations, bounds):
    population = generate_initial_population(population_size, bounds)
    for generation in range(num_generations):
        fitness = evaluate_fitness(population)
        parents = selection(population, fitness, population_size // 2)
        offspring = crossover(parents, crossover_rate)
        offspring = mutation(offspring, mutation_rate, bounds)
        population = offspring + parents
    best_solution = min(population, key=objective_function)
    return best_solution

# Set parameters
population_size = 100
mutation_rate = 0.1
crossover_rate = 0.5
num_generations = 100
bounds = (-10, 10)

# Run the Genetic Algorithm
best_solution = genetic_algorithm(population_size, mutation_rate, crossover_rate, num_generations,
bounds)
print("Best solution:", best_solution)

```


Program 2

Particle Swarm Optimization for Function Optimization

Algorithm:

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07/11/2020

Lab: Particle Swarm Optimization.

→ PSO is an optimization algorithm inspired by the social behaviour of animals such as bird flocking or fish schooling. The goal is to find optimal or near-optimal solution to complex problems by having multiple 'particles' explore the solution space. Each particle adjusts its position based on its own experience and the experience of its neighbours, ultimately converging towards the best solution.

Algorithm Code Implementation

```
import numpy as np

def RastriginFunction(x):
    A = 10
    return A * len(x) + sum((x_i**2 - A * np.cos(2 * np.pi * x_i)) for x_i in x)

class Particle:
    def __init__(self, dim):
        self.position = np.random.uniform(-5.12, 5.12, dim)
        self.velocity = np.random.uniform(-1, 1, dim)
        self.best_position = np.copy(self.position)
        self.best_value = RastriginFunction(self.position)

class PSO:
    def __init__(self, num_particles, dim, inertia_weight, cognitive_coeff, social_coeff, max_iter):
        self.num_particles = num_particles
        self.dim = dim
        self.inertia_weight = inertia_weight
        self.cognitive_coeff = cognitive_coeff
        self.social_coeff = social_coeff
        self.max_iter = max_iter
        self.swarm = [Particle(dim) for _ in range(num_particles)]
        self.global_best_position = self.swarm[0].best_position
        self.global_best_value = self.swarm[0].best_value

    def optimize(self):
        for iteration in range(self.max_iter):
            for particle in self.swarm:
                fitness_value = RastriginFunction(particle.position)
                if fitness_value < particle.best_value:
                    particle.best_value = fitness_value
                    particle.best_position = np.copy(particle.position)

                if fitness_value < self.global_best_value:
                    self.global_best_value = fitness_value
                    self.global_best_position = np.copy(particle.best_position)

            for particle in self.swarm:
                r1 = np.random.rand(self.dim)
                r2 = np.random.rand(self.dim)

                particle.velocity = (self.inertia_weight * particle.velocity +
                                     self.cognitive_coeff * r1 * (particle.best_position - particle.position) +
                                     self.social_coeff * r2 * (self.global_best_position - particle.position))

                particle.position += particle.velocity

        return self.global_best_position, self.global_best_value
```

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```
self.cognitive_coeff = cognitive_coeff
self.social_coeff = social_coeff
self.max_iter = max_iter
self.swarm = [Particle(dim) for _ in range(num_particles)]
self.global_best_position = self.swarm[0].best_position
self.global_best_value = self.swarm[0].best_value

def optimize(self):
    for iteration in range(self.max_iter):
        for particle in self.swarm:
            fitness_value = RastriginFunction(particle.position)
            if fitness_value < particle.best_value:
                particle.best_value = fitness_value
                particle.best_position = np.copy(particle.position)

            if fitness_value < self.global_best_value:
                self.global_best_value = fitness_value
                self.global_best_position = np.copy(particle.best_position)

        for particle in self.swarm:
            r1 = np.random.rand(self.dim)
            r2 = np.random.rand(self.dim)

            particle.velocity = (self.inertia_weight * particle.velocity +
                                 self.cognitive_coeff * r1 * (particle.best_position - particle.position) +
                                 self.social_coeff * r2 * (self.global_best_position - particle.position))

            particle.position += particle.velocity

        return self.global_best_position, self.global_best_value
```


Code:

```
import numpy as np

# Define the Rastrigin function
def rastrigin_function(x):
    A = 10
    return A * len(x) + sum(x_i**2 - A * np.cos(2 * np.pi * x_i) for x_i in x)

# Particle Swarm Optimization Algorithm
class Particle:
    def __init__(self, dim):
        self.position = np.random.uniform(-5.12, 5.12, dim) # Random position
        self.velocity = np.random.uniform(-1, 1, dim) # Random velocity
        self.best_position = np.copy(self.position) # Personal best position
        self.best_value = rastrigin_function(self.position) # Personal best value

class PSO:
    def __init__(self, num_particles, dim, inertia_weight, cognitive_coeff, social_coeff, max_iter):
        self.num_particles = num_particles
        self.dim = dim
        self.inertia_weight = inertia_weight
        self.cognitive_coeff = cognitive_coeff
        self.social_coeff = social_coeff
        self.max_iter = max_iter
        self.swarm = [Particle(dim) for _ in range(num_particles)]
        self.global_best_position = self.swarm[0].best_position
        self.global_best_value = self.swarm[0].best_value

    def optimize(self):
        for iteration in range(self.max_iter):
            for particle in self.swarm:
                # Evaluate fitness
                fitness_value = rastrigin_function(particle.position)

                # Update personal best
                if fitness_value < particle.best_value:
                    particle.best_value = fitness_value
                    particle.best_position = np.copy(particle.position)

                # Update global best
                if fitness_value < self.global_best_value:
                    self.global_best_value = fitness_value
                    self.global_best_position = np.copy(particle.best_position)

            # Update velocities and positions
            for particle in self.swarm:
                r1 = np.random.rand(self.dim)
```

```

r2 = np.random.rand(self.dim)

# Update velocity
particle.velocity = (self.inertia_weight * particle.velocity +
                    self.cognitive_coeff * r1 * (particle.best_position - particle.position) +
                    self.social_coeff * r2 * (self.global_best_position - particle.position))

# Update position
particle.position += particle.velocity

return self.global_best_position, self.global_best_value

# Parameters
num_particles = 30 # Number of particles
dim = 2 # Dimensionality of the problem
inertia_weight = 0.7 # Inertia weight
cognitive_coeff = 1.5 # Cognitive coefficient
social_coeff = 1.5 # Social coefficient
max_iter = 100 # Maximum number of iterations

# Run PSO
pso = PSO(num_particles, dim, inertia_weight, cognitive_coeff, social_coeff, max_iter)
best_position, best_value = pso.optimize()

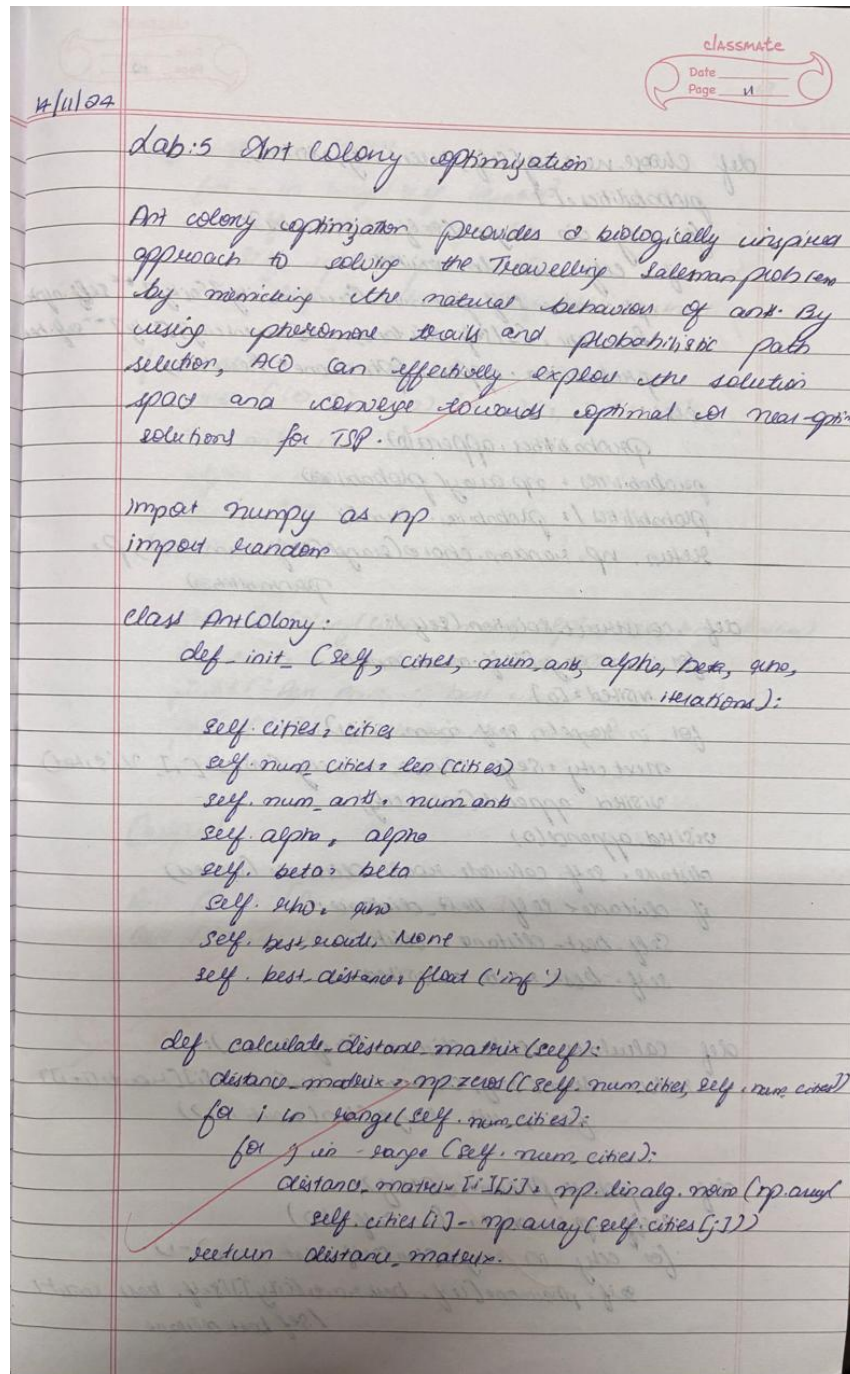
print("Best Position:", best_position)
print("Best Value:", best_value)

```

Program 3

Ant Colony Optimization for the Traveling Salesman Problem

Algorithm:



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```

def choose_next_city(self, current_city, visited):
    probabilities = []
    for city in range(self.num_cities):
        if city not in visited:
            pheromone = self.pheromone[current_city][city] ** self.alpha
            heuristic = (1 / self.distance[current_city][city]) ** self.beta
            probabilities.append(pheromone * heuristic)
        else:
            probabilities.append(0)
    probabilities = np.array(probabilities)
    probabilities /= probabilities.sum()
    return np.random.choice(range(self.num_cities), p=probabilities)

def construct_solution(self):
    for i in range(self.num_ant):
        visited = [0]
        for j in range(1, self.num_cities):
            next_city = self.choose_next_city(visited[-1], visited)
            visited.append(next_city)
        visited.append(0)
        distance = self.calculate_route_distance(visited)
        if distance < self.best_distance:
            self.best_distance = distance
            self.best_route = visited

def calculate_route_distance(self, route):
    return sum(self.distance_matrix[route[i]][route[i+1]]
                for i in range(len(route)-1))

def update_pheromone(self):
    self.pheromone *= (1 - self.rho)
    for city in range(len(self.best_route)-1):
        self.pheromone[self.best_route[city]][self.best_route[city+1]]
            / self.best_distance

```

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```

def optimize(self):
    for i in range(self.iterations):
        self.construct_solution()
        self.update_pheromone()
        return self.best_route, self.best_distance

if __name__ == '__main__':
    cities = [(0, 0), (1, 2), (2, 4), (3, 1), (4, 3)]
    num_ant = 10
    alpha = 1.0
    beta = 2.0
    rho = 0.5
    iterations = 10
    ac = AntColony(cities, num_ant, alpha, beta, rho, iterations)
    best_route, best_distance = ac.optimize()
    print("Best Route:", best_route)
    print("Best Distance:", best_distance)

```

Output:

Best Route: [0, 1, 2, 4, 3, 0]
Best Distance: 18.10654940

Code:

```
import numpy as np
import random

class AntColony:
    def __init__(self, cities, num_ants, alpha, beta, rho, iterations):
        self.cities = cities
        self.num_cities = len(cities)
        self.num_ants = num_ants
        self.alpha = alpha
        self.beta = beta
        self.rho = rho
        self.iterations = iterations
        self.pheromone = np.ones((self.num_cities, self.num_cities)) # Initial pheromone levels
        self.distance_matrix = self.calculate_distance_matrix()
        self.best_route = None
        self.best_distance = float('inf')

    def calculate_distance_matrix(self):
        distance_matrix = np.zeros((self.num_cities, self.num_cities))
        for i in range(self.num_cities):
            for j in range(self.num_cities):
                distance_matrix[i][j] = np.linalg.norm(np.array(self.cities[i]) - np.array(self.cities[j]))
        return distance_matrix

    def choose_next_city(self, current_city, visited):
        probabilities = []
        for city in range(self.num_cities):
            if city not in visited:
                pheromone = self.pheromone[current_city][city] ** self.alpha
                heuristic = (1 / self.distance_matrix[current_city][city]) ** self.beta
                probabilities.append(pheromone * heuristic)
            else:
                probabilities.append(0)
        probabilities = np.array(probabilities)
        probabilities /= probabilities.sum() # Normalize
        return np.random.choice(range(self.num_cities), p=probabilities)
```

```

def construct_solution(self):
    for _ in range(self.num_ants):
        visited = [0] # Start from the first city
        for _ in range(1, self.num_cities):
            next_city = self.choose_next_city(visited[-1], visited)
            visited.append(next_city)
        visited.append(0) # Return to the starting city
        distance = self.calculate_route_distance(visited)
        if distance < self.best_distance:
            self.best_distance = distance
            self.best_route = visited

def calculate_route_distance(self, route):
    return sum(self.distance_matrix[route[i]][route[i + 1]] for i in range(len(route) - 1))

def update_pheromones(self):
    # Evaporate pheromones
    self.pheromone *= (1 - self.rho)
    # Add pheromones based on the best route found
    for city in range(len(self.best_route) - 1):
        self.pheromone[self.best_route[city]][self.best_route[city + 1]] += 1 / self.best_distance

def optimize(self):
    for _ in range(self.iterations):
        self.construct_solution()
        self.update_pheromones()
    return self.best_route, self.best_distance

```

Example usage:

```

if __name__ == "__main__":
    # Define cities as (x, y) coordinates
    cities = [(0, 0), (1, 2), (2, 4), (3, 1), (4, 3)]

```

ACO parameters

```

num_ants = 10
alpha = 1.0 # importance of pheromone
beta = 2.0 # importance of heuristic

```



```
rho = 0.5 # pheromone evaporation rate
iterations = 10

aco = AntColony(cities, num_ants, alpha, beta, rho, iterations)
best_route, best_distance = aco.optimize()

print("Best Route:", best_route)
print("Best Distance:", best_distance)
```

Program 4

Cuckoo Search (CS)

Algorithm:

21-11-24

Lab 6: Cuckoo Search

Cuckoo search is a ^{new} inspired optimization algorithm based on the brood parasitism of some cuckoo species. This behaviour involves laying eggs in the nest of other birds, leading to the optimization of search strategies. CS uses Lévy flights to generate new solutions, promoting global search capabilities and avoiding local minima.

```

import numpy as np
import random
from scipy.special import gamma

def levy_flight(beta=1.5, alpha=1):
    sign = 1 if np.power(gamma(1+beta)) * np.sin(np.pi * (beta/2) / gamma((1+beta)/2)) > 0 else -1
    p1 = beta/2
    p2 = (beta/2)**2
    u = np.random.normal(0, sigma=1, size=1)
    v = np.random.normal(0, 1, size=1)
    step = u / (np.power(np.abs(v), 1/beta))
    return sign * step

def initialize_population(n_nests, n_dim, lower_bound, upper_bound):
    return np.random.uniform(lower_bound, upper_bound, (n_nests, n_dim))

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

def cuckoo_search(n_nests, n_dim, lower_bound, upper_bound, max_iter, p0=0.25):
    nests = initialize_population(n_nests, n_dim, lower_bound, upper_bound)
    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 iteration in range(max_iter):
        for i in range(n_nests):
            step = levy_flight(alpha, n_dim)
            new_nest = nests[i] + step * (nests[i] - best_nest)
            new_nest = np.clip(new_nest, lower_bound, upper_bound)
            new_fitness = fitness_function(new_nest)
            if new_fitness < fitness[i]:
                nests[i] = new_nest
                fitness[i] = new_fitness
            if new_fitness < best_fitness:
                best_nest = new_nest
                best_fitness = new_fitness

        for i in range(n_nests):
            if random.random() < p0:
                nests[i] = np.random.uniform(lower_bound, upper_bound, n_dim)
                fitness[i] = fitness_function(nests[i])

        if (iteration+1) * 100 == 0 or iteration == max_iter-1:
            print(f"Iteration {iteration+1}, Best Fitness: {best_fitness}")
            return best_nest, best_fitness

n_nests = 25
n_dim = 10
lower_bound = -5
upper_bound = 5
    
```

Code:

```
import numpy as np
import random
from scipy.special import gamma # Import gamma from scipy.special

# Levy flight function
def levy_flight(beta=1.5, d=1):
    sigma_u = np.power((gamma(1 + beta) * np.sin(np.pi * beta / 2) / gamma((1 + beta) / 2) * beta *
                        np.cos(np.pi * beta / 2) ** 2), 1 / beta)
    u = np.random.normal(0, sigma_u, size=d)
    v = np.random.normal(0, 1, size=d)
    step = u / np.power(np.abs(v), 1 / beta)
    return step

# Initialize population (nests)
def initialize_population(n_nests, n_dim, lower_bound, upper_bound):
    return np.random.uniform(lower_bound, upper_bound, (n_nests, n_dim))

# Fitness function (Example: Sphere function)
def fitness_function(x):
    return np.sum(x ** 2)

# Cuckoo Search Algorithm
def cuckoo_search(n_nests, n_dim, lower_bound, upper_bound, max_iter, pa=0.25):
    # Step 1: Initialize nests randomly
    nests = initialize_population(n_nests, n_dim, lower_bound, upper_bound)

    # Step 2: Evaluate fitness of all nests
    fitness = np.array([fitness_function(nest) for nest in nests])

    # Track the best solution found so far
    best_idx = np.argmin(fitness)
    best_nest = nests[best_idx]
    best_fitness = fitness[best_idx]

    # Start iterations
    for iteration in range(max_iter):
        # Generate new solutions using Levy flight
        for i in range(n_nests):
            # Generate a new solution by Levy flight
            step = levy_flight(d=n_dim)
            new_nest = nests[i] + step * (nests[i] - best_nest)

            # Apply boundary conditions
            new_nest = np.clip(new_nest, lower_bound, upper_bound)

            # Evaluate the new solution
```

```

new_fitness = fitness_function(new_nest)

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

# Update the best solution if necessary
if new_fitness < best_fitness:
    best_nest = new_nest
    best_fitness = new_fitness

# Abandon some of the worst nests and generate new random solutions
for i in range(n_nests):
    if random.random() < pa: # with probability pa
        nests[i] = np.random.uniform(lower_bound, upper_bound, n_dim)
        fitness[i] = fitness_function(nests[i])

# Print progress every 100 iterations
if (iteration + 1) % 100 == 0 or iteration == max_iter - 1:
    print(f"Iteration {iteration + 1}, Best Fitness: {best_fitness}")

return best_nest, best_fitness

# Parameters
n_nests = 25 # Number of nests
n_dim = 10 # Dimensionality of the problem
lower_bound = -5 # Lower bound for the search space
upper_bound = 5 # Upper bound for the search space
max_iter = 1000 # Maximum number of iterations
pa = 0.25 # Probability of abandoning the worst nests

# Run Cuckoo Search
best_solution, best_solution_fitness = cuckoo_search(n_nests, n_dim, lower_bound, upper_bound,
max_iter, pa)

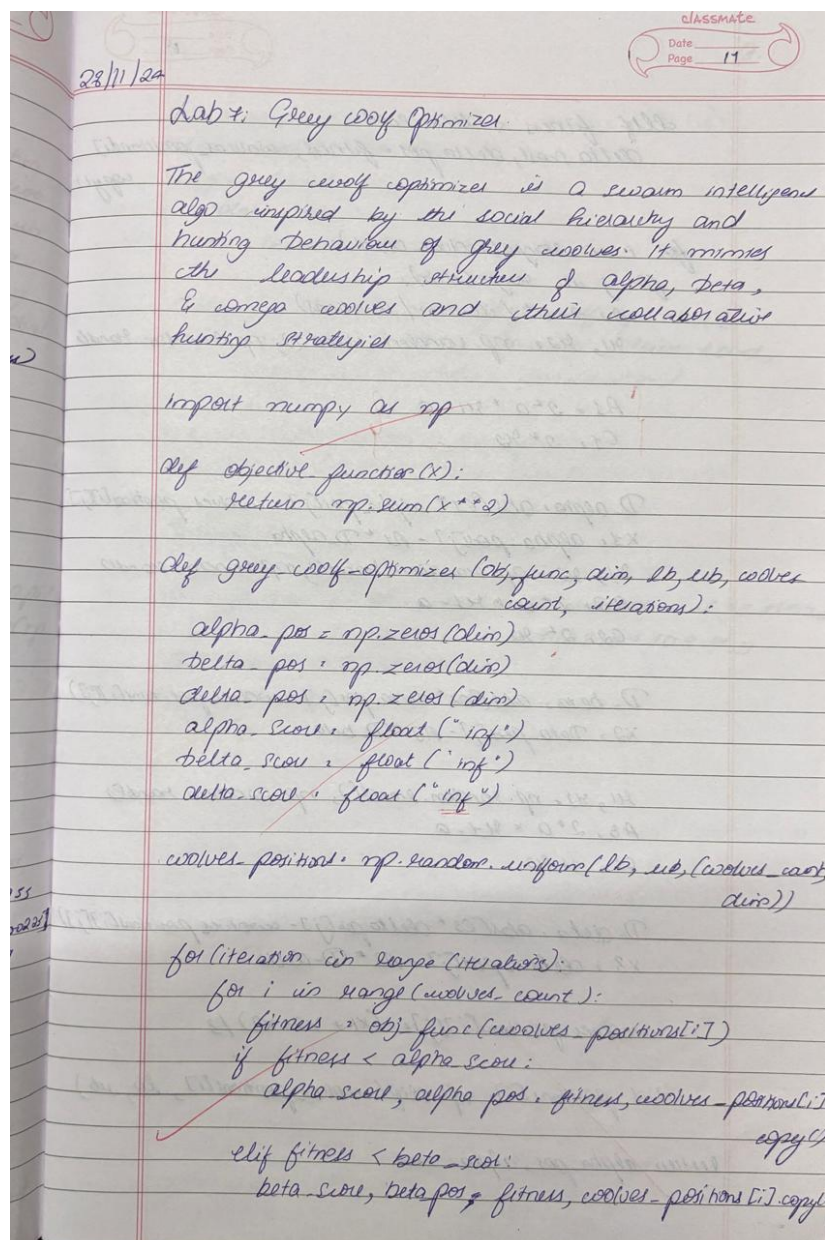
print("\nBest solution found: ", best_solution)
print("Best fitness value: ", best_solution_fitness)

```

Program 5

Grey Wolf Optimizer (GWO)

Algorithm:



clip_fitness < delta_pos:

delta_pos, delta_pos = fitness, evolves_position[i]

for i in range(evolve_count):

for j in range(dim):

a = a + iterations / iterations

x1, x2 = np.random.rand(1), np.random.rand(1)

A1 = 2 * a * x1 - a

C1 = 2 * x2

D_alpha = abs(C1 * alpha_pos[j] - evolves_position[i][j])

x1 = alpha_pos[j] - A1 * D_alpha

x1, x2 = np.random.rand(1), np.random.rand(1)

A2 = 2 * a * x1 - a

C2 = 2 * x2

D_beta = abs(C2 * beta_pos[j] - evolves_position[i][j])

x2 = beta_pos[j] - A2 * D_beta

x1, x2 = np.random.rand(1), np.random.rand(1)

A3 = 2 * a * x1 - a

C3 = 2 * x2

D_delta = abs(C3 * delta_pos[j] - evolves_position[i][j])

x3 = delta_pos[j] - A3 * D_delta

evolves_position[i][j] = (x1 + x2 + x3) / 3

evolve_position[i] = np.clip(evolve_position[i], lb, ub)

return alpha_pos, alpha_pos

dimension = 5

lower_bound = -10

upper_bound = 10

evolves = 50

max_iteration = 50

best_position, test_score = grey_wolf_optimizer(
objective_function, dimension, lower_bound,
upper_bound, evolves, max_iteration)

print("Best position", best_position)

print("Best score", test_score)

Output

Best position: 16.39416682 -6.4005065 -6.5951078

7.98219731 -7.16874410E-07

Best Score: 0.4045629534

~~0.11~~

Code:

```
import numpy as np

# Define the objective function to be minimized
def objective_function(x):
    return np.sum(x**2) # Example: Sphere function (minimization problem)

# Grey Wolf Optimizer (GWO) implementation
def grey_wolf_optimizer(obj_func, dim, lb, ub, wolves_count, iterations):
    # Initialize alpha, beta, and delta wolves' positions
    alpha_pos = np.zeros(dim)
    beta_pos = np.zeros(dim)
    delta_pos = np.zeros(dim)

    # Initialize alpha, beta, and delta wolves' fitness values
    alpha_score = float("inf") # Best fitness
    beta_score = float("inf") # Second-best fitness
    delta_score = float("inf") # Third-best fitness

    # Initialize the positions of all wolves
    wolves_positions = np.random.uniform(lb, ub, (wolves_count, dim))

    # Main loop for optimization
    for iteration in range(iterations):
        for i in range(wolves_count):
            # Calculate the fitness of the current wolf
            fitness = obj_func(wolves_positions[i])

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

        # Update the positions of wolves
        for i in range(wolves_count):
            for j in range(dim):
                # Coefficients
                a = 2 - 2 * (iteration / iterations) # Linearly decreases from 2 to 0
                r1, r2 = np.random.rand(), np.random.rand()

                A1 = 2 * a * r1 - a
                C1 = 2 * r2
```

```

D_alpha = abs(C1 * alpha_pos[j] - wolves_positions[i][j])
X1 = alpha_pos[j] - A1 * D_alpha

r1, r2 = np.random.rand(), np.random.rand()
A2 = 2 * a * r1 - a
C2 = 2 * r2

D_beta = abs(C2 * beta_pos[j] - wolves_positions[i][j])
X2 = beta_pos[j] - A2 * D_beta

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

D_delta = abs(C3 * delta_pos[j] - wolves_positions[i][j])
X3 = delta_pos[j] - A3 * D_delta

# Update position
wolves_positions[i][j] = (X1 + X2 + X3) / 3

# Enforce bounds
wolves_positions[i] = np.clip(wolves_positions[i], lb, ub)

# Return the best solution
return alpha_pos, alpha_score

# Parameters
dimension = 5 # Number of variables
lower_bound = -10 # Lower bound of variables
upper_bound = 10 # Upper bound of variables
wolves = 30 # Number of wolves in the pack
max_iterations = 50 # Maximum number of iterations

# Run the GWO algorithm
best_position, best_score = grey_wolf_optimizer(
    objective_function, dimension, lower_bound, upper_bound, wolves, max_iterations
)

print("Best position:", best_position)
print("Best score:", best_score)

```

Program 6

Parallel Cellular Algorithms and Programs

Algorithm:

19/12/29

Lab 8: Parallel Cellular Algorithms

Parallel cellular algorithms are inspired by the functioning of biological cells that operate in a highly parallel and distributed manner.

```
import numpy as np
def roaster-function(x):
    A=10
    return A * len(x) + sum((xi**2 - A * np.cos(2 * np.pi * xi))
    for xi in x)

class ParallelCellularAlgorithm:
    def __init__(self, grid_size=(5,5), dim=2, neighborhood='moor', iterations=100):
        self.grid_size = grid_size
        self.dim = dim
        self.neighborhood = neighborhood
        self.bounds = bounds
        self.grid = np.random.uniform(bounds[0], bounds[1],
        grid_size[0], grid_size[1], dim)
        self.fitness = np.zeros([grid_size[0], grid_size[1]])

    def evaluate_fitness(self):
        for i in range(self.grid_size[0]):
            for j in range(self.grid_size[1]):
                self.fitness[i,j] = roaster_func(self.grid[i,j])

    def get_neighbors(self, i, j):
        neighbors = []
        directions = []
        if self.neighborhood == 'moor':
            directions = [(-1,-1), (-1,0), (-1,1)]
        elif self.neighborhood == 'von_neuman':
            directions = [(-1,0), (0,-1), (0,1)]

    def update_status(self):
        new_grid = self.grid.copy()
        for i in range(self.grid_size[0]):
            for j in range(self.grid_size[1]):
                neighbors = self.get_neighbors(i,j)
                self.grid[i,j] = new_grid[i,j]

    def run(self):
        for iteration in range(self.iterations):
            self.evaluate_fitness()
            self.update_status()
            print(f"Iteration {iteration+1} / {self.iterations}")

        print("Optimization Complete")
        print(f"Best Solution: {self.best_solution}")
        print(f"Best Fitness: {self.best_fitness}")

if __name__ == "__main__":
    pca = ParallelCellularAlgorithm(grid_size=(10,10),
    dim=2, neighborhood='moor')
    pca.run()
```

Output.

Optimization Complete
Best Solution: [-0.0337945 -0.0340272]
Best Fitness: 0.451717

Code:

```
import numpy as np

def rastrigin_function(x):
    """
    Rastrigin function for optimization. It has a global minimum at x = 0.
     $f(x) = 10n + \sum(x_i^2 - 10 * \cos(2 * \pi * x_i))$ 
    """
    A = 10
    return A * len(x) + sum((xi ** 2 - A * np.cos(2 * np.pi * xi)) for xi in x)

class ParallelCellularAlgorithm:
    def __init__(self, grid_size=(5, 5), dim=2, neighborhood='moore', iterations=100, bounds=(-5.12, 5.12)):
        """
        Initialize the parameters for the cellular algorithm.

        :param grid_size: Tuple defining the dimensions of the grid (rows, columns)
        :param dim: Dimensionality of the solution space
        :param neighborhood: Type of neighborhood ('moore' or 'von_neumann')
        :param iterations: Number of iterations to run
        :param bounds: Lower and upper bounds for the solution space
        """
        self.grid_size = grid_size
        self.dim = dim
        self.neighborhood = neighborhood
        self.iterations = iterations
        self.bounds = bounds

        # Initialize the grid with random solutions
        self.grid = np.random.uniform(bounds[0], bounds[1], size=(grid_size[0], grid_size[1], dim))
        self.fitness = np.zeros((grid_size[0], grid_size[1]))

        # Store the best solution and its fitness
        self.best_solution = None
        self.best_fitness = float('inf')

    def evaluate_fitness(self):
        """
        Evaluate fitness of each cell in the grid.
        """
        for i in range(self.grid_size[0]):
            for j in range(self.grid_size[1]):
                self.fitness[i, j] = rastrigin_function(self.grid[i, j])

        # Update the best solution
        if self.fitness[i, j] < self.best_fitness:
```

```

        self.best_fitness = self.fitness[i, j]
        self.best_solution = self.grid[i, j].copy()

def get_neighbors(self, i, j):
    """
    Get neighboring cells based on the chosen neighborhood type.

    :param i: Row index of the current cell
    :param j: Column index of the current cell
    :return: List of neighboring cells
    """
    neighbors = []
    directions = []
    if self.neighborhood == 'moore':
        # Moore neighborhood: 8 neighbors (including diagonals)
        directions = [(-1, -1), (-1, 0), (-1, 1), (0, -1), (0, 1), (1, -1), (1, 0), (1, 1)]
    elif self.neighborhood == 'von_neumann':
        # Von Neumann neighborhood: 4 neighbors (up, down, left, right)
        directions = [(-1, 0), (0, -1), (0, 1), (1, 0)]

    for di, dj in directions:
        ni, nj = i + di, j + dj
        if 0 <= ni < self.grid_size[0] and 0 <= nj < self.grid_size[1]:
            neighbors.append(self.grid[ni, nj])
    return neighbors

def update_states(self):
    """
    Update the state of each cell based on its neighbors.
    """
    new_grid = self.grid.copy()
    for i in range(self.grid_size[0]):
        for j in range(self.grid_size[1]):
            neighbors = self.get_neighbors(i, j)

            # Calculate the average position of neighbors
            avg_neighbor = np.mean(neighbors, axis=0)

            # Move the current cell slightly towards the average neighbor position
            new_grid[i, j] = self.grid[i, j] + 0.1 * (avg_neighbor - self.grid[i, j])

            # Ensure the new solution stays within bounds
            new_grid[i, j] = np.clip(new_grid[i, j], self.bounds[0], self.bounds[1])

    self.grid = new_grid

def run(self):
    """

```

Run the Parallel Cellular Algorithm.

```
"""
```

```
for iteration in range(self.iterations):
```

```
    # Step 1: Evaluate fitness of the current grid
```

```
    self.evaluate_fitness()
```

```
    # Step 2: Update the states of all cells
```

```
    self.update_states()
```

```
    # Print progress
```

```
    print(f"Iteration {iteration + 1}/{self.iterations}, Best Fitness: {self.best_fitness:.6f}")
```

```
print("Optimization complete.")
```

```
print(f"Best Solution: {self.best_solution}")
```

```
print(f"Best Fitness: {self.best_fitness:.6f}")
```

```
if __name__ == "__main__":
```

```
    # Initialize and run the algorithm
```

```
    pca = ParallelCellularAlgorithm(grid_size=(10, 10), dim=2, neighborhood='moore', iterations=50)
```

```
    pca.run()
```


Program 7

Optimization via Gene Expression Algorithms

Algorithm:

19/11/20

Lab 9: Optimization via Gene Expression Algorithm.

Gene Expression Algo are inspired by the biological process of gene expression in living organisms. The process involves the translation of genetic information encoded in DNA into functional proteins.

import numpy as np
import random

def maximum-function(x):
A = 10
return $A \cdot \ln(x) + \sum_{i=1}^n (x_i + 2 - A \rightarrow \text{np.cos}(2 \cdot \text{np.pi} \cdot x_i))$ for x_i in x

class GeneExpressionAlgorithm:
def __init__(self, pop_size=100, gene_len=10, bounds=(-5.12, 5.12),
self.pop_size, pop_size
self.gene_len, gene_len
self.mutation_rate, mutation_rate
self.bounds, bounds.

def random_gene_sequence(self):
return np.random.uniform(self.bounds[0], self.bounds[1],
size=self.gene_len)

def evaluate_fitness(self, gene_sequence):
return maximum-function(gene_sequence)

def selection(self):
tournament_size = 3
selected = []
for i in range(self.pop_size):
tournament = random.sample(self.population, tournament_size)
return selected.

def mutate(self, gene_sequence):
for i in range(len(gene_sequence)): if random.random() < self.mutation_rate:
gene_sequence[i] += np.random.uniform(-1.0, 1.0)
gene_sequence[i] = np.clip(gene_sequence[i], self.bounds[i],
return gene_sequence.

def gene_expression(self, gene_sequence):
return gene_sequence.

def run(self):
for generation in range(self.generations):
fitness_values = [self.evaluate_fitness(ind) for ind in self.population]
best_idx = np.argmax(fitness_values)
selected_pop = self.selection()
next_population = []
for i in range(0, self.pop_size, 2):
parent1 = selected_pop[i]
parent2 = selected_pop[i+1]
offspring1, offspring2 = self.crossover(p1, p2)
next_pop.append(self.mutate(offspring1,))
self.population = next_population

Print ("Optimization complete")
Print ("Best Sol: ", self.best_sol)
Print ("Best Fitness: ", self.best_fitness)

gla.run()

D/P Optimization complete
Best Sol: [-0.02796841 -0.01908592]
Best Fitness: 0.226974.

Code:

```
import numpy as np
import random
```

```
def rastrigin_function(x):
```

```
    """
```

```
    Rastrigin function for optimization. It has a global minimum at x = 0.
```

```
     $f(x) = 10n + \sum(x_i^2 - 10 * \cos(2 * \pi * x_i))$ 
```

```
    """
```

```
    A = 10
```

```
    return A * len(x) + sum((xi ** 2 - A * np.cos(2 * np.pi * xi)) for xi in x)
```

```
class GeneExpressionAlgorithm:
```

```
    def __init__(self, population_size=100, gene_length=10, mutation_rate=0.01, crossover_rate=0.7,
generations=100, bounds=(-5.12, 5.12)):
```

```
        """
```

```
        Initialize the parameters for the Gene Expression Algorithm.
```

```
        :param population_size: Number of genetic sequences in the population
```

```
        :param gene_length: Length of each genetic sequence
```

```
        :param mutation_rate: Probability of mutation
```

```
        :param crossover_rate: Probability of crossover
```

```
        :param generations: Number of generations to evolve
```

```
        :param bounds: Lower and upper bounds for the solution space
```

```
        """
```

```
        self.population_size = population_size
```

```
        self.gene_length = gene_length
```

```
        self.mutation_rate = mutation_rate
```

```
        self.crossover_rate = crossover_rate
```

```
        self.generations = generations
```

```
        self.bounds = bounds
```

```
        # Initialize population with random solutions
```

```
        self.population = [self.random_gene_sequence() for _ in range(population_size)]
```

```
        self.best_solution = None
```

```
        self.best_fitness = float('inf')
```

```
    def random_gene_sequence(self):
```

```
        """Generate a random gene sequence within bounds."""
```

```
        return np.random.uniform(self.bounds[0], self.bounds[1], size=self.gene_length)
```

```
    def evaluate_fitness(self, gene_sequence):
```

```
        """Evaluate the fitness of a gene sequence using the Rastrigin function."""
```

```
        return rastrigin_function(gene_sequence)
```

```
    def selection(self):
```

```
        """Select individuals for reproduction using tournament selection."""
```

```
        tournament_size = 3
```

```
        selected = []
```

```
25 for _ in range(self.population_size):
```

```
        tournament = random.sample(self.population, tournament_size)
```

```
        tournament_fitness = [self.evaluate_fitness(ind) for ind in tournament]
```

```

        winner = tournament[np.argmin(tournament_fitness)]
        selected.append(winner)
    return selected

def crossover(self, parent1, parent2):
    """Perform crossover between two parents to produce offspring."""
    if random.random() < self.crossover_rate:
        point = random.randint(1, self.gene_length - 1)
        offspring1 = np.concatenate((parent1[:point], parent2[point:]))
        offspring2 = np.concatenate((parent2[:point], parent1[point:]))
        return offspring1, offspring2
    return parent1, parent2

def mutate(self, gene_sequence):
    """Apply mutation to a gene sequence."""
    for i in range(len(gene_sequence)):
        if random.random() < self.mutation_rate:
            gene_sequence[i] += np.random.uniform(-1.0, 1.0)
            gene_sequence[i] = np.clip(gene_sequence[i], self.bounds[0], self.bounds[1])
    return gene_sequence

def gene_expression(self, gene_sequence):
    """Translate a genetic sequence into a functional solution (no changes needed here)."""
    return gene_sequence

def run(self):
    """Run the Gene Expression Algorithm."""
    for generation in range(self.generations):
        # Step 1: Evaluate fitness and find the best solution
        fitness_values = [self.evaluate_fitness(ind) for ind in self.population]
        best_idx = np.argmin(fitness_values)
        if fitness_values[best_idx] < self.best_fitness:
            self.best_fitness = fitness_values[best_idx]
            self.best_solution = self.population[best_idx].copy()

        # Step 2: Selection
        selected_population = self.selection()

        # Step 3: Crossover and Mutation
        next_population = []
        for i in range(0, self.population_size, 2):
            parent1 = selected_population[i]
            parent2 = selected_population[i + 1]
            offspring1, offspring2 = self.crossover(parent1, parent2)
            next_population.append(self.mutate(offspring1))
            next_population.append(self.mutate(offspring2))

        self.population = next_population

```

```
print("Optimization complete.")
print(f"Best Solution: {self.best_solution}")
print(f"Best Fitness: {self.best_fitness:.6f}")

if __name__ == "__main__":
    # Initialize and run the Gene Expression Algorithm
    gea = GeneExpressionAlgorithm(population_size=50, gene_length=2, generations=50)
    gea.run()
```