

## LAB. 1

Summarising optimisation problems:

### 1. Genetic Algorithm

It simulates the process of natural selection. A random population  $P$  is selected and a string of random characters represent chromosomes. Fitness of population is determined, selection of the fittest is performed then crossover & mutation is performed until the best fitness level is reached.

### 2. Particle Swarm Optimisation

It is inspired by the social behaviour of bird flocking or fish schooling.

Each particle with the swarm evaluates its position with the help of a fitness function, adjusting their velocity and position based on best position and converge on most optimal solution.

### 3. Ant Colony Optimisation

It is based on the idea of foraging of ants, which is used to find solutions for the complex travelling salesman problem.

It executes by mimicking the behaviour of how ants choose the shortest path between their food sources and their nest.

## Introduction to Genetic Algorithms

- Historical context :

In the early 1950s, experiments in AI and evolutionary computation, with Nils Bärtoli creating the first software mimicking reproduction and mutation.

Genetic algorithms were developed by John Holland in the 1960s inspired by the principles of natural selection & evolution (Darwin's theory).

- Abstract

GAs are powerful optimisation tools inspired by the processes of natural evolution (selection, mutation & crossover). By simulating the process of natural selection, GAs search for optimal solutions.

- Working Process (with terminologies) :

In a GA, the fitness function first evaluates how well each chromosome performs in solving the problem.

- Then chromosomes, which are potential solutions encoded as arrays, pass through solution process.
- Selection favours the fitter chromosomes for reproduction, increasing the likelihood of producing better offspring.
- Crossover combines two selected parent chromosomes.

Mutation then introduces random changes to some chromosomes to achieve genetic diversity & preventing premature convergence.

These steps are performed until optimal solution is reached.

We usually perform maximising or minimising of a particular function. There is a fitness value  $f(x)$  associated to every  $x$  value & a selection probability.

Selection is performed w.r.t. the function to find the optimal solution & further crossover & mutation of these 5 bit string are performed to generate new fitness population.

Date - 16/10/24

## PROGRAM - 1

Program - Implement a Genetic algorithm using python to solve basic optimisation problem.

import numpy as np

def func(x):

    return x\*\*2

Population-size = 30

num-generation = 50

mutation-rate = 0.1

crossover-rate = 0.1

range-min, range-max = -10, 10

def initialise\_population(size, minval, maxval):

    return np.random.uniform(minval, maxval, size)

def evaluate\_fitness(population):

    return np.array([func(n) for n in population])

def select\_parents(population, fitness):

    total\_fitness = np.sum(fitness)

    selection\_probs = fitness / total\_fitness

    return np.random.choice(population, size=2)

P = selection\_probs

def crossover(Parent1, Parent2):

    if (np.random.rand() < crossover\_rate):

        return (Parent1 + Parent2) / 2

    return Parent1

```

def mutate(offspring, min-val, max-val):
    if np.random.rand() < mutation_rate:
        return np.random.uniform(min-val, max-val)
    return offspring

def genetic_algorithm():
    population = initialise_population(Population_size,
                                         range_min, range_max)
    best_solution = None
    best_fitness = -inf

    new_population = []
    for _ in range(Population_size):
        parent_1, parent_2 = select_parents(population, fitness)
        offspring = crossover(parent_1, parent_2)
        offspring = mutate(offspring, range_min, range_max)
        new_population.append(offspring)

    population = np.array(new_population)

    return best_solution, best_fitness

```

Get best\_n, best\_value = genetic\_algorithm()

~~Print ("best\_n : {best\_n}, Max. value of f(x) : {best\_value}")~~

Output:

best n : - 9.98

Max value of f(n) : 99.611

Date: 23/10/24

## PROGRAM - 2

### Particle Swarm Optimisation

The technique is inspired by the flocking behaviour of birds, swarming of insects and schooling of fish. It is a heuristic search algorithm to find optimal solution to a problem.

#### Algorithm:

1. **Initialisation**: Define the problem domain & initialize the population of particles randomly.
2. **Evaluation**: Evaluate the fitness of each particle (based on its position)
3. **Update the personal best**: Update the personal best position of each particle.
4. **Update the global best position** based on the best of all particles.
5. **Update velocity and position** of each particle based on its current position and velocity, and the personal & global best positions.
6. **Termination**: Repeat steps 2-5 for a fixed number of iterations or until a stopping criterion is met i.e target fitness value is achieved.

```
import numpy as np  
def objective_function(n):  
    return np.sum(n**2)  
  
class particle:  
    def __init__(self, bounds):  
        self.position = np.array([np.random.uniform(bound[0], bound[1]) for bound in bounds])  
        self.velocity = np.zeros_like(self.position)  
        self.best_pos = self.position.copy()  
        self.best_value = float('inf')
```

```
def update_personal_best(self, value):
```

```
    if value < self.best_value:
```

 ~~self.best\_value = value~~ ~~self.best\_pos = self.best\_position.copy()~~

```
def PSO(objective_function, bounds, num_particles=30,  
        mom_iterations=20, inertia_weight=0.7,  
        cognitive_coefficient=1.5, social_coefficient=1.5):
```

```
    particles = [Particle(bounds) for _ in range(num_particles)]
```

```
    global_best_position = None
```

```
    global_best_value = float('inf')
```

```
    for particle in particles:
```

```
        fitness_value = objective_function(particle.position)
```

```
        particle.update_personal_best(fitness_value)
```

If fitness\_value < global\_best\_value

global\_best\_value = fitness\_value

global\_best\_position = particle.position.copy()

for particle in particles:

    Inertia-component = inertia\_weight \* particle\_velocity

    Cognitive-component = cognitive\_coeff \* np.random.random()  
                          \* (particle.best\_position - particle.position)

    Social-component = social\_coeff \* np.random.random() +  
                          (global\_best\_position - particle.position)

    particle.velocity = Inertia-component + Cognitive-component  
                          + Social-component

    particle.position += particle.velocity

for i in range(len(bounds)):

    particle.position[i] = np.clip(particle.position[i],  
                                bounds[i][0], bounds[i][1])

    print(f"Iteration {iteration + 1} / {max\_iterations}

    Global-best\_value : {global\_best\_value})")

return global\_best\_position, global\_best\_value

bounds = [(-5, 5), (-5, 5)]

best\_position, best\_value = pso(objective\_function, bounds)

print(f"Best Position : {best\_position}")  
print(f"Best Value : {best\_value}")

Output:

Iteration 1/20, Global Best Value: 1.0462397424940557

Best Position → [-0.00476255 -0.001899]

Best Value : 2.6288113052694566 e-05

Sept  
23/10/24

## PROGRAM - 3

### Ant Colony Optimisation

It is a probabilistic technique for solving computational problems that can be reduced to finding good paths through graphs.

### Algorithm:

Initialise necessary parameters

Generate ant population

Find the fitness value associated with each ant

Find best solution through selection methods

Update pheromone trail

### Code:

```
import numpy as np
```

```
import random
```

```
def create_distance_matrix(n_cities):
```

```
    np.random.seed(0)
```

```
    matrix = np.random.randint(1, 100, size=(n_cities, n_cities))
```

```
    np.fill_diagonal(matrix, 0)
```

```
    return matrix
```

n\_cities = 10

n\_ants = 20

n\_iterations = 50

alpha = 1

beta = 2

evaporation\_rate = 0.5

initial\_pheromone = 1

distance-matrix = create-distance-matrix ( $n$ -cities)  
pheromone-matrix = np.ones (( $n$ -cities,  $n$ -cities)) \* initial-  
Pheromone

class Ant:

def \_\_init\_\_(self, n-cities):

self.n-cities = n-cities

self.route = []

self.distance\_travelled = 0

def select-next-city (self, current-city, visited):

probabilities = []

for city in range (self.n-cities):

if city not in visited:

pheromone = pheromone-matrix[current-city][city]  
\*\* alpha

heuristic = (1 / distance-matrix [current-city][city])  
\*\* beta

probabilities.append (pheromone \* heuristic)

else

probabilities.append (0)

Probabilities = np.array (probabilities) / sum (probabilities)

next-city = np.random.choice (range (self.n-cities),  
p=probabilities)

return next-city

def find\_route (self):

current-city = random.randint (0, self.n-cities - 1)

self.route = [current-city]

visited = set (self.route)

Self-distance-travelled += distance matrix [current-city][next-city]

def update-pheromone(ants):

global pheromone-matrix

pheromone-matrix \*= (1 - evaporation-rate)

for ant in ants:

for i in range (len(ant.route)-1):

city-from = ant.route[i]

city-to = ant.route[i+1]

pheromone-matrix [city-from][city-to] += 1.0 /  
ant.distance-travelled

pheromone-matrix [city-to][city-from] += 1.0 /  
ant.distance-travelled

def ant\_colony-optimisation():

best-route = None

best-distance = float('inf')

~~for~~ for iteration in range (n-iterations):

ants = [Ant(n-cities) for \_ in range (n-ants)]

for ant in ants:

ant.find\_route()

if ant.distance-travelled < best-distance:

best-distance = ant.distance-travelled

best-route = ant.route

Update-pheromones  
print(f"Iteration {iteration+1}: Best distance =  
{best\_distance}")

return best-route, best-distance

best-route, best-distance = ant-colony-optimisation()

print(f"Best route found : {best\_route} with distance.  
{best\_distance}")

### Output:

Iteration 1: Best distance = 187

Iteration 2: Best distance = 130

Iteration 3: Best distance = 130

Iteration 50: Best distance = 118.

Best route found = [7, 4, 2, 6, 1, 3, 8, 9, 0, 5, 7]  
with distance : 118.

Date: 13/11/24

## PROGRAM-4

### Cuckoo Search Algorithm.

This algorithm is based on the brood parasitic behaviour of cuckoo species.

#### Usage:

It is used to optimise solutions in cloud computing, data mining, software, IoT and pattern recognition.

#### Code:

```
import numpy as np
```

```
def levy_flight (beta, dim):
```

$$\text{sigma\_u} = (\text{np.math.gamma}(1+\beta) * \text{np.sin}(\text{np.pi} * \beta / 2) / (\text{np.math.gamma}((1+\beta)/2) * \beta^{1+\beta}))^{1/\beta}$$

```
u = np.random.normal (0, sigma_u, dim)
```

```
v = np.random.normal (0, 1, dim)
```

```
step = u / np.abs(v) ** (1/beta)
```

```
return step
```

```
def objective_function (x) :
```

```
return np.sum(x ** 2)
```

```
def cuckoo_search (obj_func, dim, population_size,  
max_iter, alpha=0.01, beta=1.5,  
pa=0.25) :
```

nest = np.random.uniform(-10, 10, (population\_size, dim))

fitness = np.apply\_along\_axis(obj\_func, 1, nests)

best\_nest\_idn = np.argmax(fitness)

best\_nest = nests[best\_nest\_idn]

best\_fitness = fitness[best\_nest\_idn]

for iteration in range(max\_iter):

for i in range(population\_size):

step = levy\_flight(beta, dim)

new\_nest = nests[i] + alpha \* step

new\_nest = np.clip(new\_nest, -10, 10)

new\_fitness = obj\_func(new\_nest)

if new\_fitness < fitness[i]:

nests[i] = new\_nest

fitness[i] = new\_fitness

for i in range(population\_size):

if np.random.rand() < p\_a:

nests[i] = np.random.uniform(-10, 10, dim)

current\_best\_idn = np.argmax(fitness)

current\_best\_fitness = fitness[current\_best\_idn]

if current\_best\_fitness < best\_fitness:

best\_fitness = current\_best\_fitness

best\_nest = nests [current\_best\_idx]

print (f"Iteration {iteration+1}/{max\_iter}, Best fitness  
{best\_fitness}")

return best\_nest, best\_fitness

#parameters

dim = 2

Population\_size = 20

max\_iter = 100

alpha = 0.01

beta = 1.5

Pa = 0.25

best\_solution, best\_value = cuckoo\_search (objective function,  
dim, population\_size, max\_iter, alpha, beta, pa)

print (f"Best solution found: {best\_solution}")

print (f"Objective function value: {best\_value}")

### Output:

Iteration 1/100, Best Fitness: 14.158607731707642

Iteration 2/100, Best Fitness: 1.829912696458187

Iteration 3/100, Best Fitness: 1.829912696458187

Iteration 100/100, Best Fitness: 0.1786865018893165

Best Solution found: [3.77880199 -3.5320835]

# PROGRAM - 5

## GREY WOLF OPTIMISER

This algorithm is inspired by the social behaviour and hunting technique of grey wolves. The purpose of this algorithm is to find the optimal solution by simulating the social structure and hunting tactics of grey wolves in herd.

Usage : It is used in engineering design problems for structural optimisation and mechanical design.

In machine learning , it is used in feature selection and training neural networks.

Code :

```
import numpy as np
def fitness_function (Position):
    return np.sum(position ** 2)
```

```
def grey_wolf_optimisation (fitness_function, dim, n_wolves,
    max_iter, lower_bound, upper_bound):
```

```
    wolves = np.random.uniform (lowerbound, upperbound,
        (n_wolves, dim))
```

```
    fitness = np.apply_along_axis (fitness_function, 1, wolves)
```

```
    alpha_pos = wolves [np.argmin (fitness)]
```

```
    alpha_score = np.min (fitness)
```

```
    fitness [np.argmin(fitness)] = float ("inf")
```

```
    beta_pos = wolves [np.argmin(fitness)]
```

```
    beta_score = np.min (fitness)
```

`fitness [np.argmax(fitness)] = float ("inf")`

`delta_pos = wolves [np.argmax(fitness)]`

`delta_score = np.mean(fitness)`

for t in range (n\_wolves):

$r_1, r_2 = \text{np.random.random(dim)}, \text{np.random.random(dim)}$

$A1, C1 = 2 * a * r_1 - a, 2 * r_2$

$r_1, r_2 = \text{np.random.random(dim)}, \text{np.random.random(dim)}$

$A2, C2 = 2 * a * r_1 - a, 2 * r_2$

$r_1, r_2 = \text{np.random.random(dim)}, \text{np.random.random(dim)}$

$A3, C3 = 2 * a * r_1 - a, 2 * r_2$

$D_{alpha} = \text{np.abs}(C1 * alpha_pos - wolves[i])$

$D_{beta} = \text{np.abs}(C2 * beta_pos - wolves[i])$

$D_{delta} = \text{np.abs}(C3 * delta_pos - wolves[i])$

$X1 = alpha_pos - A1 * D_{alpha}$

$X2 = beta_pos - A2 * D_{beta}$

$X3 = delta_pos - A3 * D_{delta}$

$wolves[i] = (X1 + X2 + X3) / 3$

`wolves = np.clip(wolves, lower_bound, upper_bound)`

`fitness = np.apply_along_axis(fitness_function, 1, wolves)`

`alpha_idn = np.argmax(fitness)`

if `fitness[alpha_idn] < alpha_score`:

`alpha_pos = wolves[alpha_idn]`

`alpha_score = fitness[alpha_idn]`

$\text{beta\_idn} = \text{wp.argmax(fitness)}$

if  $\text{fitness}[\text{beta\_idn}] < \text{beta\_score}$  and  $\text{fitness}[\text{beta\_idn}] > \text{alpha\_score}$ :

$\text{beta\_pos} = \text{wolves}[\text{beta\_idn}]$

$\text{beta\_score} = \text{fitness}[\text{beta\_idn}]$

$\text{delta\_idn} = \text{wp.argmax(fitness)}$

if  $\text{fitness}[\text{delta\_idn}] < \text{delta\_score}$  and  $\text{fitness}[\text{delta\_idn}] > \text{beta\_score}$ :

$\text{delta\_pos} = \text{wolves}[\text{delta\_idn}]$

$\text{delta\_score} = \text{fitness}[\text{delta\_idn}]$

`printf ("Iteration {++1}/{monitors}, Best Fitness :  
{alpha-score}")`

`return alpha_pos, alpha-score.`

`dim = 5`

`n_wolves = 30`

`max_iter = 100`

`lower_bound = -10`

`upper_bound = 10`

`best_position, best_score = grey_wolf_optimisation(fitness_function,  
dim, n_wolves, max_iter, lower_bound, upper_bound)`

`print ("Best Position:", best_position)`

`print ("Best Score:", best_score.)`

**Output :**

`Best Position : [-0.53923125 0.6760124 -1.847277247  
-3.96605915 -0.30903728]`

`Best Score : 1.1558289302571314`

## PROGRAM - 6

### Parallel Cellular Algorithm

It is designed to solve complex problem that can be broken down into local, independent task executed.

The application of this algorithm are in image processing, physical simulation, biological simulation, optimisation and machine learning.

#### Algorithm:

- 1) Initialisation : Define a grid  $G$  of size  $n \times m$ , where each cell  $G(x,y)$  has initial state from a set  $S$

$$S = \{0, 1\} \text{ for binary states}$$

$$G = \{G(x,y) \mid 0 \leq x < n, 0 \leq y < m\}$$

- 2) Define neighbourhood : Define neighbourhood  $N(G(x,y))$  for each cell  $G(x,y)$ . Common choice, include Moore neighbourhood  $\rightarrow$  all 8 surrounding cells  $3 \times 3$  grid including the center.

- 3) Update Rule : Each cell new state  $G'(x,y)$  is complete based on its current state and the state of its neighbour using an update function.

$$G'(x,y) = f(G(x,y), \{G(x',y') \mid (x',y') \in N(G(x,y))\})$$

#### Example updates rule

- i) Parallel Update : The update are computed for all cells  $(x,y)$  in parallel

$$G'(x,y) = f(G(x,y), N(G(x,y)) \# (x,y))$$

- 5) Duration: Repeat the update iteration for  $T$  time step or until convergence
- 6) Termination: Stop the algorithm when the predefined no. of iteration is reached or when grid stabilizes.

### Iteration 1

$$\begin{bmatrix} 0 & 0 & 1 & 0 & 1 \\ 0 & 0 & 0 & 0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 0 & 1 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 0 & 1 & 0 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 0 & 0 & 0 & 1 \end{bmatrix}$$

### Iteration 2

$$\begin{bmatrix} 0 & 1 & 1 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 0 & 1 & 0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 1 & 1 & 0 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 1 & 1 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 1 & 0 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 1 & 1 & 1 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 1 & 0 & 1 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 1 & 1 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 1 & 0 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 1 & 1 & 1 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 1 & 0 & 1 & 1 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 1 & 1 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 0 & 1 & 0 & 1 & 0 \end{bmatrix}$$

$$\begin{bmatrix} 1 & 1 & 1 & 1 & 1 \end{bmatrix}$$

## PROGRAM - 7

### Gene Expression Algorithm

This is an evolutionary algorithm that draws attention from the biological process of gene expression. Its primary purpose is to optimise complex problems and devise solutions computational challenges by simulating the mechanisms of evolution and gene transcription of living organic.

Usage: Used in optimisation problems, data mining and machine learning, symbolic regression; control systems, genetic programming, bioinformatics

### Algorithm:

- 1) Generate an initial population  $P(0)$  of  $N$  chromosomes (individuals). Each chromosome  $C_i$  is represented as a string of symbols (genotype) ( $i = \{g_1, g_2, \dots, g_L\}$ ,  $L = \text{length of chromosome}$ )

$g_i$  is a gene selected from a predefined terminal:

$$T = \{n_1, n_2, n_3, \dots, n_n\} \quad F = \{+, -, \times, \div\}$$

- 2) Decode the chromosome: Convert the chromosome  $C_i$  into a phenotype using a decoding process.

$$f(x) = n_1 + (n_2 \times n_3) \dots$$

- 3) Fitness Evaluation: Evaluate the fitness  $F(c_i)$  of each chromosome using a problem specific

fitness function.

$$F(C) = \frac{1}{n} \sum_{j=1}^n (y_j - f(x_j))^2$$

$y_j$  = actual output for input  $x_j$

$f(x_j)$  = Predicted output for  $c_i$

minimize the fitness function.

4. Selection : Select the top-performing chromosomes for a population based on their fitness  $F(i)$ . Use methods like Roulette wheel selection or tournament selection.

5. Genetic Operations :

i) Crossover:

$$O_1, O_2 = \text{Crossover}(c_1, c_2)$$

ii) Mutation:

$$c'_i = \text{Mutate}(c_i)$$

iii) Transportation : Move parts of the new chromosomes to a new location within the chromosome.

6. Replacement : Replace the old population  $P(f)$  with new population  $P(t(t))$  based on fitness :

$$P(t_f) = \{O_1, O_2, \dots, O_n\}$$

7. Termination : Repeat steps 2-6 for a no. of generations  $G$  or until termination criterion is met.

Initial Population:

## Generation 2:

## Chromosome

O : 2509.7331

1:2657.1015

2:14f

3 : 2509.7331

4 : 2546.5782

Final:

## Fitness.

## Chromosome

O: 2582.4173

1 : 2582.4173

2: 3147.2067

3:1289,2067

4 : 1416. 6804

Best Solution:

Chromosome Fitness: 1259.2067

$$\{ \varrho_0, \dots, \varrho_n, 0 \} = \{ \varrho \}$$