

EXERCISE 1

The goal of simulated annealing is to maximize the performance of a model relative to a fitness function defined over the model parameters you are seeking to optimize. In Exercise 2, you will apply simulated annealing coded in scipy to a one-dimensional problem in which you seek the value of X that maximizes the output of a function called `fitnessFunction`. But first you must write code that implements the function, which is the product of m and n : $m = 1 + \cos(0.04v)^2$ $n = \exp(-v^2 / (20000))$

Entrée [2]: `import math`

```
def fitnessFunction(v):
    m = 1 + math.cos(0.04*v)**2
    n = math.exp(-v**2 / (20000))
    return m*n
```

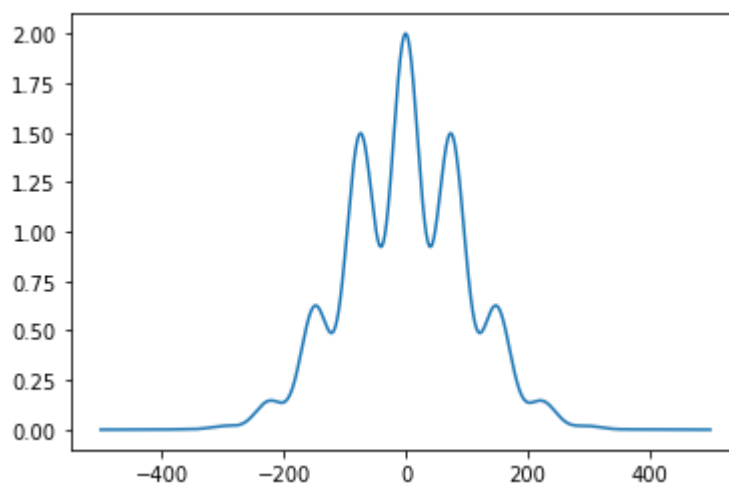
EXERCISE 2

Display the fitness surface. To do this, make the wave fitness with 1000 points starting at $x = -500$ and ending at $x = 500$ (using Data > Change Wave Scaling). Set the y values in fitness to the corresponding value of the `fitnessFunction()`. Display fitness and include the graph in your lab report.

Entrée [4]: `import matplotlib.pyplot as plt`
`import numpy as np`

```
def graph_fitness():
    v = np.arange(-500., 500., .5)
    fitness = []
    for x in v:
        fitness.append(fitnessFunction(x))
    plt.plot(v, fitness)

graph_fitness()
```



EXERCISE 3

Run the algorithm with an initial X value of 250 and an initial temperature of 10. Plot the trajectory of the annealing algorithm along the fitness surface, Include the graph in your lab report. Slowly increase the temperature until you find the peak about 90% of the time. Report this temperature and explain its significance in terms of the relationship between the search radius and the dimensions of the width of the fitness curve.

```

Entrée [38]: ▶ import random
import math
import matplotlib.pyplot as plt

debug=False
interval = (-500, 500)

def f(x):
    # penalized version of fitnessFunction to avoid going out of interval
    ro = 1
    result = fitnessFunction(x)
    if x < min(interval):
        result -= ro * (min(interval)-x)**2
    elif x > max(interval):
        result -= ro * (x-max(interval))**2
    return result

def perturbation(fraction=1):
    # small move within interval length
    amplitude = (max(interval) - min(interval)) * fraction / 10
    delta = amplitude * random.uniform(-1, 1)
    return delta

def simulated_annealing(init_state, t0, tend=.01, alpha=0.9, max_nit=20):

    global states
    global costs
    global iterations

    iterations=[]
    costs = []
    states=[]
    nit=0

    current_state = init_state
    current_e = f(current_state)
    t = t0

    while (t > tend):
        nit += 1
        next_state = current_state + perturbation(0.5)
        next_e = f(next_state)
        delta_e = next_e - current_e
        delta_e = - delta_e
        if delta_e < 0:
            current_state = next_state
        else:
            if math.exp(-delta_e / t) > random.random():
                current_state = next_state
            current_e = f(current_state)
        t = t * alpha
        iterations.append(nit)
        costs.append(fitnessFunction(current_state))
        states.append(current_state)

    return current_state

```

```
global states
global costs

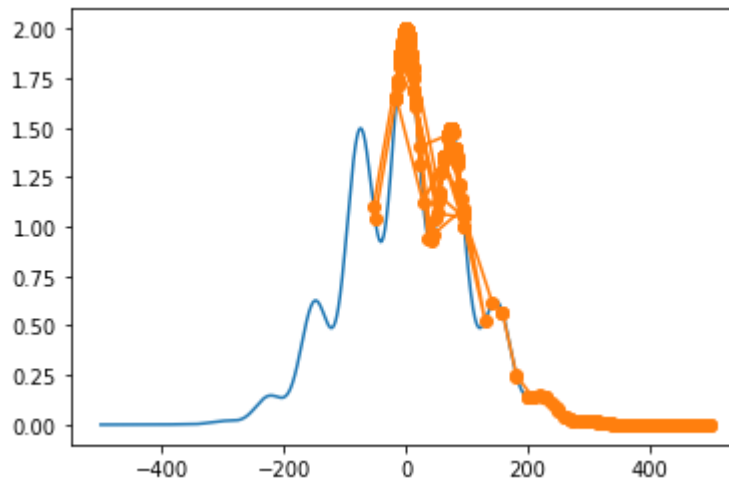
init_value = 250
init_temp = 10
final_temp = 0.001
alpha = 0.99

result = simulated_annealing(init_value, t0=init_temp, tend=final_temp, alpha)
print("Maximal value after annealing: %.4f found at x=%.4f after %d iterations")

graph_fitness()
plt.plot(states, costs, "-o")
```

Maximal value after annealing: 2.0000 found at x=0.1202 after 917 iterations

Out[38]: [<matplotlib.lines.Line2D at 0x19bae8fb860>]

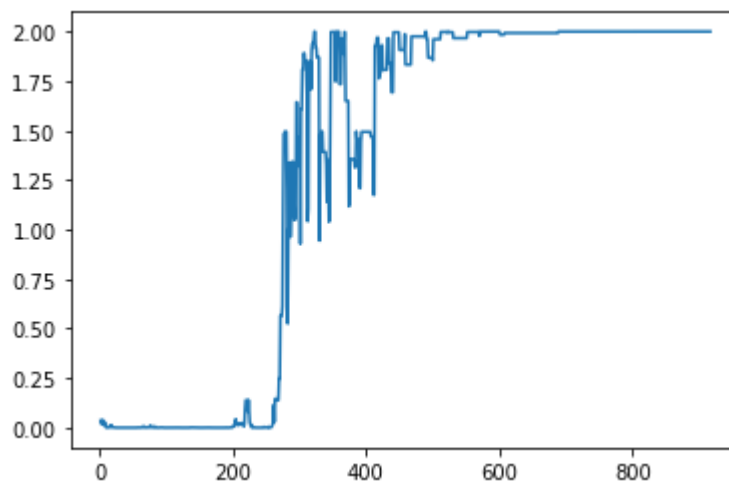


EXERCISE 4

Another informative graph is the plot of all the intermediate best fitnesses versus iteration number. Make this plot. Include the graph in your lab report

```
Entrée [39]: plt.plot(iterations, costs)
```

```
Out[39]: [<matplotlib.lines.Line2D at 0x19baea308d0>]
```



EXERCISE 5

Complete the table below by running simulation 20 times with each of 10 different values of temp shown. Plot number of failures vs temperature and number of iterations to reach the peak vs temperature. Describe and explain any trends you find.

```

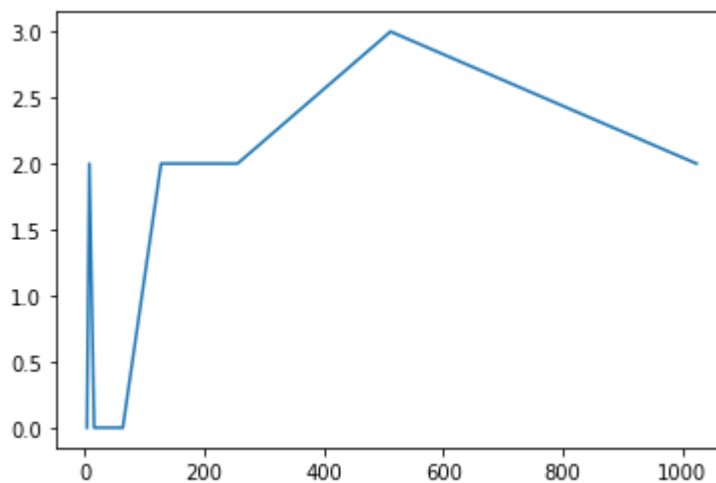
Entrée [17]: newt = 2
temperatures = []
failures = []
for i in range(0,9):
    newt *= 2
    temperatures.append(newt)
    failures.append(0)

    for j in range(0,19):
        result = simulated_annealing(init_value, t0=newt, tend=final_temp, a=0.99)
        if abs(result)>5:
            failures[i] += 1
            #print("Maximal value after annealing: %.4f found at x=%.4f after %d iterations" % (result, x, i))

plt.plot(temperatures, failures)

```

Out[17]: [<matplotlib.lines.Line2D at 0x19babd5eb70>]



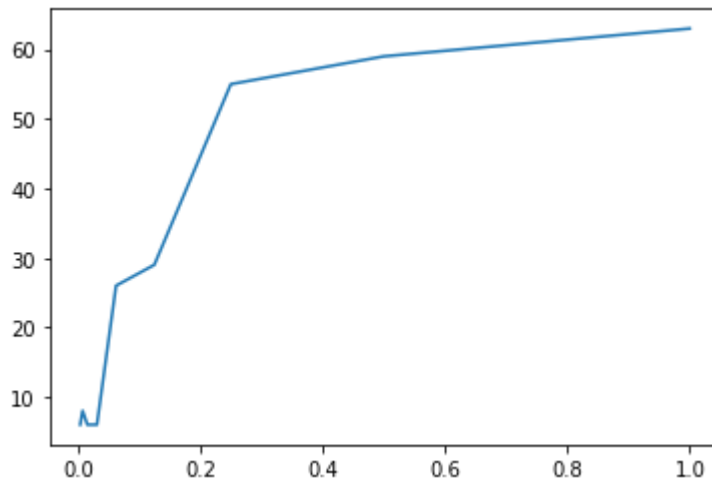
Interesting: nb of failure vs final temerature
 in this case simulated annealing seems sensitive to end temperature.
 as ΔE is very small if final temperature is too high $\exp(\Delta E/t)$ is
 always 0.99 and accept most of perturbations

```
Entrée [25]: ▶ newt = 2
temperatures = []
failures = []
for i in range(0,9):
    newt /= 2
    temperatures.append(newt)
    failures.append(0)

    for j in range(0,99):
        result = simulated_annealing(init_value, t0=10, tend=newt, alpha=alpha)
        if abs(result)>5:
            failures[i] += 1
            #print("Maximal value after annealing: %.4f found at x=%.4f after %d" % (result, x, j))

plt.plot(temperatures,failures)
```

Out[25]: [<matplotlib.lines.Line2D at 0x19bad4bda58>]



Entrée []: ▶