

While working on Fourier Series or some other Mathematical Problem, you might sometime have to work with Periodic Functions.

Periodic Functions are those that give the same value after a particular period.

So we will use this definition to define a periodic function in PYTHON.

Let's say that there is a function  $f(x)$  which is defined in the interval  $[li,lf]$  and is periodic with a period of  $T=lf-li$ .

Then the function should have the same value at:  $f(x)$ ,  $f(x+T)$ ,  $f(x+2*T)$ , ....

i.e.  $f(x)=f(x+T)=f(x+2*T)=\dots\dots$  since period= $T$ .

But I said that the function is defined only in the interval  $[li,lf]$ . So how is the computer supposed to calculate its value at  $x>lf$ ?

That's easy. Since the value of the function at  $f(x+T)$  is simply  $f(x)$ , therefore we can generalize that whenever  $x>lf$ : then,

$f(x)=f(x-T)$ . Note: We have to keep taking  $x$  back by  $T$  i.e  $(x-T)$  until it lies within  $[li,lf]$  where the function is well-defined.

Similarly what about the value of function at  $x$  less than  $(li)$  cause the function is not defined for values less than  $(li)$ ?

Again, this time we use:  $f(x)=f(x+T)$ . Note: We keep translating  $x$  forward by  $T$  i.e  $(x+T)$  until it lies within  $[li,lf]$  where the function is well-defined.

Using the above two arguments we can create a function which will make any given function defined within  $[li,lf]$  and with a period  $T$  a periodic function.

Here is the code for that:

```
# Function that will convert any given function 'f' defined in a given range '[li,lf]'
to a periodic function of period 'lf-li'
def periodicf(li,lf,f,x):
    if x>=li and x<=lf :
        return f(x)
    elif x>lf:
        x_new=x-(lf-li)
        return periodicf(li,lf,f,x_new)
    elif x<(li):
        x_new=x+(lf-li)
        return periodicf(li,lf,f,x_new)
```

In the above code I have created a function 'periodicf' which takes as arguments the limits  $[li,lf]$  for which the function 'f' (third argument) is originally defined and the fourth argument is the value of  $x$  at which I want the value of 'f'.

The above code assumes the function to be defined within  $[li,lf]$  therefore the function's starting point is  $(li)$ . However, if you want to create a function that is defined in a different manner then you will have to define 'f' correspondingly.

Following is a program which you can use to define various periodic functions such as sawtooth wave, cycloids, square wave, and triangular wave.

```
import numpy as np
import matplotlib.pyplot as plt
from matplotlib.pyplot import *
```

```

fig = figure(figsize=(8, 8), dpi=120)

# Function that will convert any given function 'f' defined in a given range '[li,lf]'
to a periodic function of period 'lf-li'
def periodicf(li,lf,f,x):
    if x>=li and x<=lf :
        return f(x)
    elif x>lf:
        x_new=x-(lf-li)
        return periodicf(li,lf,f,x_new)
    elif x<(li):
        x_new=x+(lf-li)
        return periodicf(li,lf,f,x_new)

# The periodic version of sawtooth function
def sawtoothP(li,lf,x):
    return periodicf(li,lf,sawtooth,x)

# Non-periodic sawtooth function defined for a range [-l,l]
def sawtooth(x):
    return x

# The periodic version of square function
def squareP(li,lf,x):
    return periodicf(li,lf,square,x)

# Non-periodic square wave function defined for a range [-l,l]
def square(x):
    if x>0:
        return 5
    else:
        return 0

# The periodic version of triangle function
def triangleP(li,lf,x):
    return periodicf(li,lf,triangle,x)

# Non-periodic triangle wave function defined for a range [-l,l]
def triangle(x):
    if x>0:
        return x
    else:
        return -x

# The periodic version of cycloid function
def cycloidP(li,lf,x):
    return periodicf(li,lf,cycloid,x)

# Non-periodic cycloid wave function defined for a range [-l,l]
def cycloid(x):

```

```

return np.sqrt(5**2-x**2)

if __name__ == "__main__":

    # plt.style.use('dark_background')
    plt.style.use('seaborn')
    plt.title('Periodic functions\nCycloid')

    li = -5
    lf = 5
    step_size = 0.05

    x_l = -20
    x_u = 50

    x = np.arange(x_l,x_u,step_size)
    y1 = [sawtoothP(li,lf,xi) for xi in x]
    y2 = [squareP(li,lf,xi) for xi in x]
    y3 = [triangleP(li,lf,xi) for xi in x]
    y4 = [cycloidP(li,lf,xi) for xi in x]

    x_plot = []
    y_plot1 = []
    y_plot2 = []
    y_plot3 = []
    y_plot4 = []

    x_l_plot = x_l - 15
    x_u_plot = x_l_plot + 20
    plt.xlim(x_l_plot,x_u_plot)
    plt.ylim(-6,7)

    for i in range(x.size):
        x_plot.append(x[i])
        y_plot1.append(y1[i])
        y_plot2.append(y2[i])
        y_plot3.append(y3[i])
        y_plot4.append(y4[i])
        #Sawtooth
        plt.plot(x_plot,y_plot1,c='darkkhaki')
        #Square
        plt.plot(x_plot,y_plot2,c='tomato')
        #Triangular
        plt.plot(x_plot,y_plot3,c='orange')
        #Cycloid
        plt.plot(x_plot,y_plot4,c='slateblue')
        x_l_plot = x_l_plot + step_size
        x_u_plot = x_u_plot + step_size

```

```
plt.xlim(x_l_plot,x_u_plot)
plt.pause(0.1)
plt.show()
```

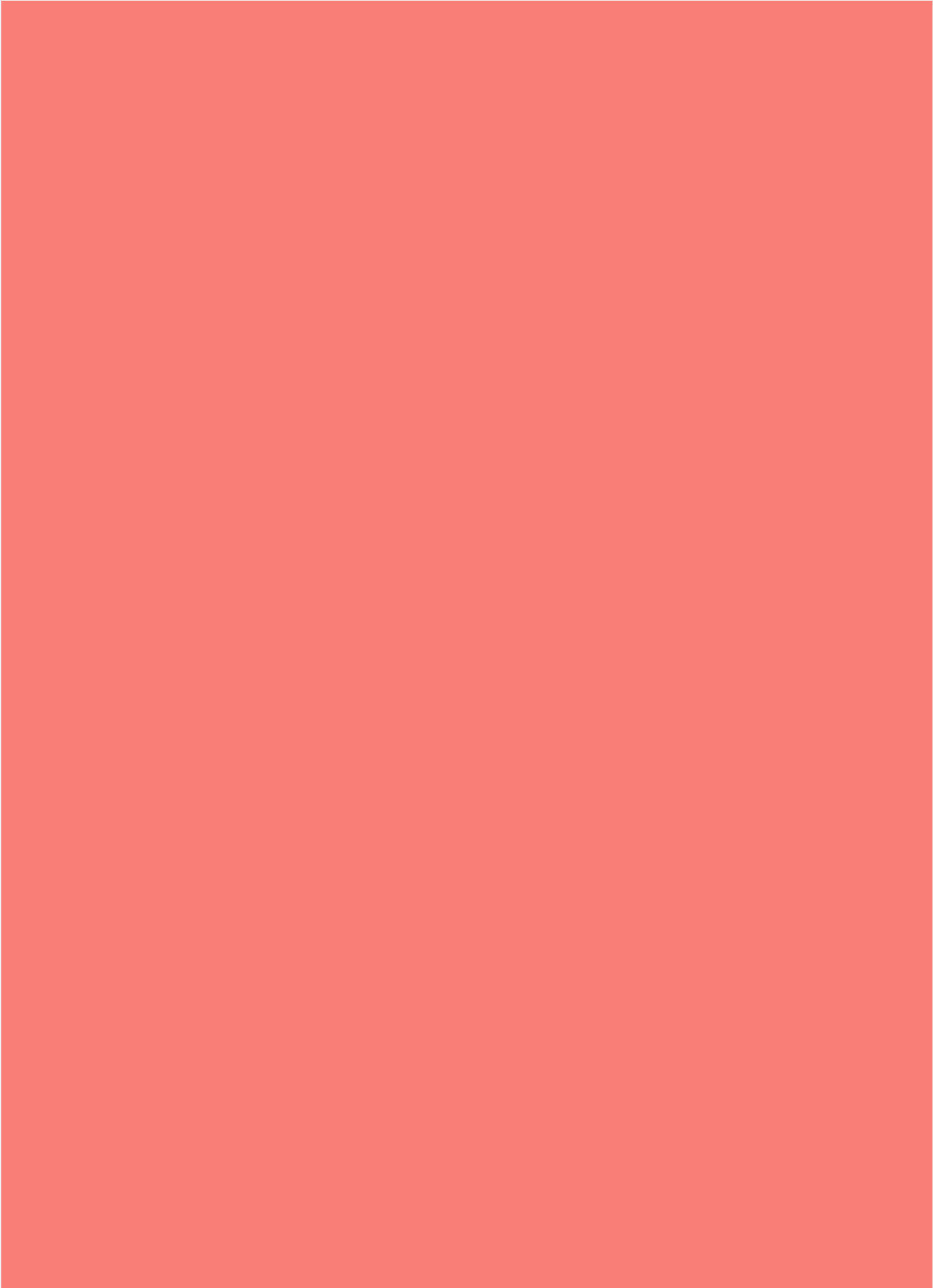
## OUTPUT

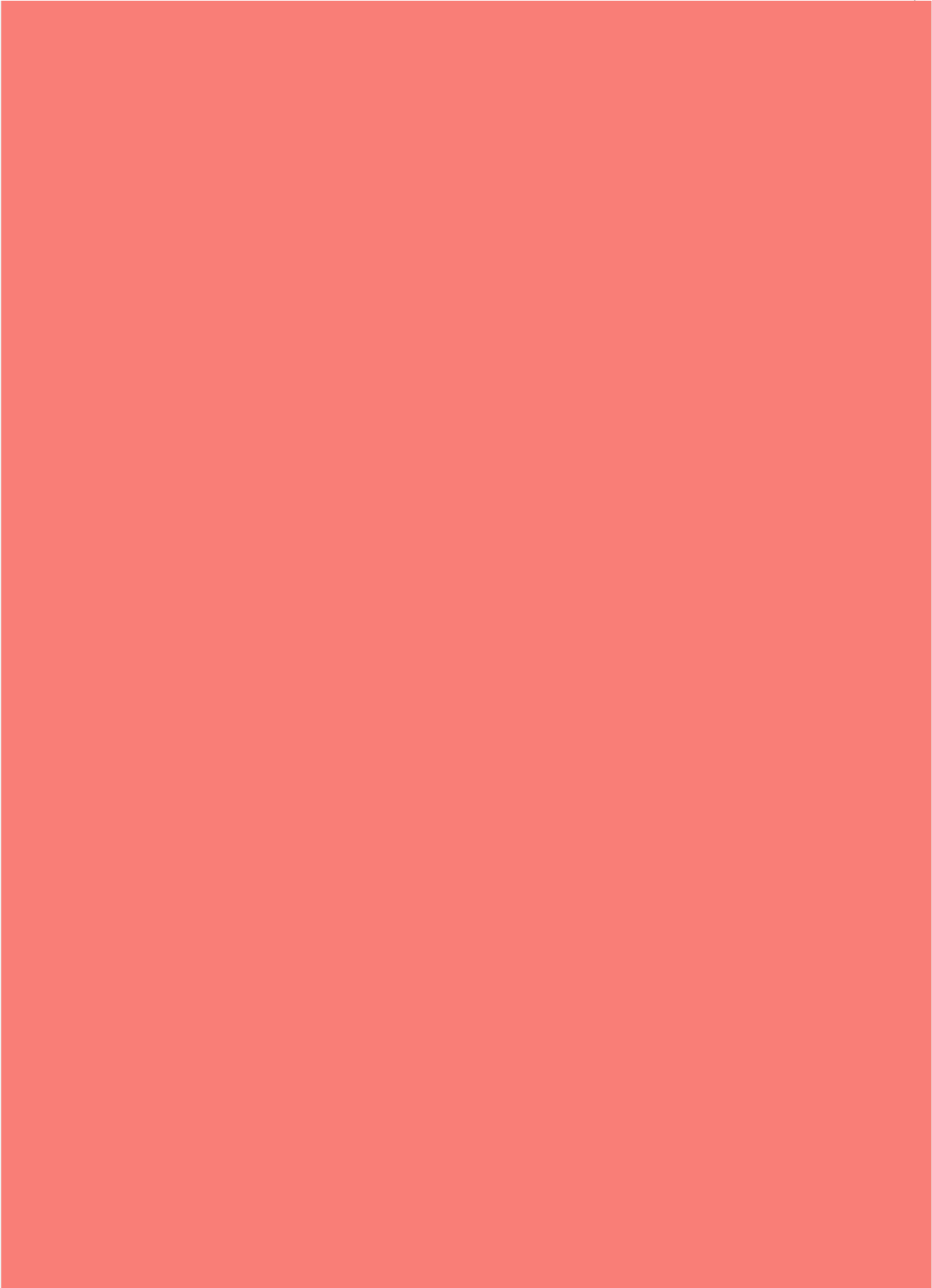


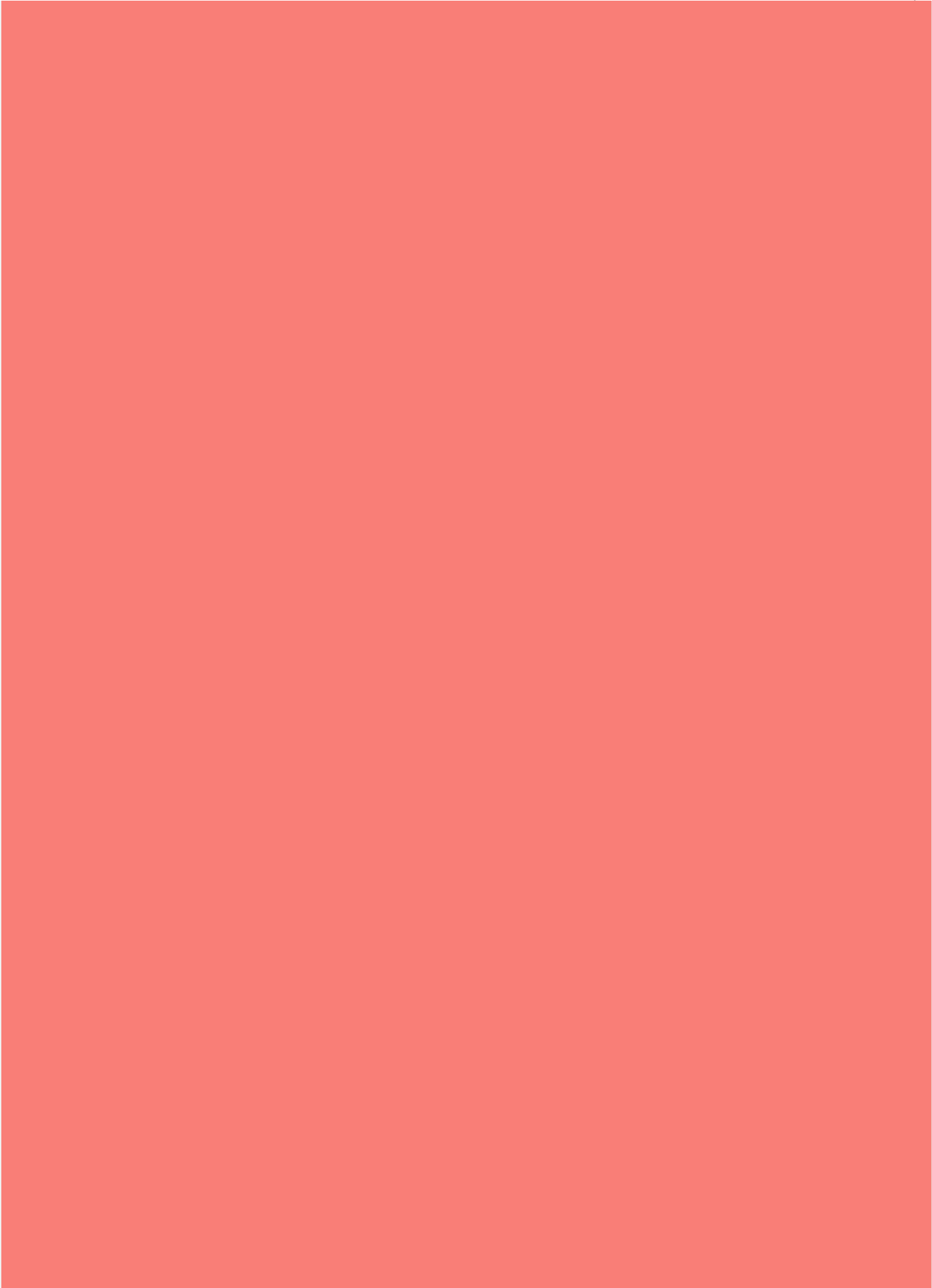
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I'm a physicist specializing in computational material science with a PhD in Physics from Friedrich-Schiller University Jena, Germany. I write efficient codes for simulating light-matter interactions at atomic scales. I like to develop Physics, DFT, and Machine Learning related apps and software from time to time. Can code in most of the popular languages. I like to share my knowledge in Physics and applications using this Blog and a YouTube channel.

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