

Homework 7

Problem 1

Vandermonde function:

a)

```
In [1]: import numpy as np
import matplotlib.pyplot as plt
import interp as lag
```

①

a)

$$V = \begin{bmatrix} 1 & x_0 & x_0^2 & \dots & x_0^{n-1} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 1 & x_1 & x_1^2 & \dots & x_1^{n-1} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 1 & x_n & x_n^2 & \dots & x_n^{n-1} \end{bmatrix}$$

so $V\vec{c} = \vec{y}$

gives $\begin{bmatrix} 1 & x_0 & x_0^2 & \dots & x_0^{n-1} & c_1 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 1 & x_1 & x_1^2 & \dots & x_1^{n-1} & c_1 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 1 & x_n & x_n^2 & \dots & x_n^{n-1} & c_n \end{bmatrix} = \begin{bmatrix} y_1 \\ \vdots \\ y_n \end{bmatrix}$

```
In [2]: def vander(V,y):
sol = np.linalg.solve(V,y)
return sol
```

b)

```
In [3]: N1=2
i1 = np.linspace(1,N1,N1)
x1= -1+(i1-1)*(2/(N1-1))
x1=x1.reshape((2,1))
y1= 1/(1+(10*x1)**2)
def V1(x,N):
v=np.ones((N,1))
#print(v)
for n in range(1,N):
#print(n)
#print(x**n)

v=np.hstack(([v,(x**n)]))
#print(v)
```

```

    return v
def polyCoefficients(x, coeffs):

    order = len(coeffs)

    y = 0
    for i in range(order):
        y += coeffs[i]*x**i
    return y

```

Testing

In [4]: `print(x1**3)`

```

[[ -1.]
 [  1.]]

```

In [5]: `print(V1(x1,N1))`
`print(y1)`
`print((x1**3))`

```

[[  1. -1.]
 [  1.  1.]]
[[0.00990099]
 [0.00990099]]
[[ -1.]
 [  1.]]

```

In [6]: `poly1 = vander(V1(x1,N1),y1)`

In [7]: `print(poly1)`

```

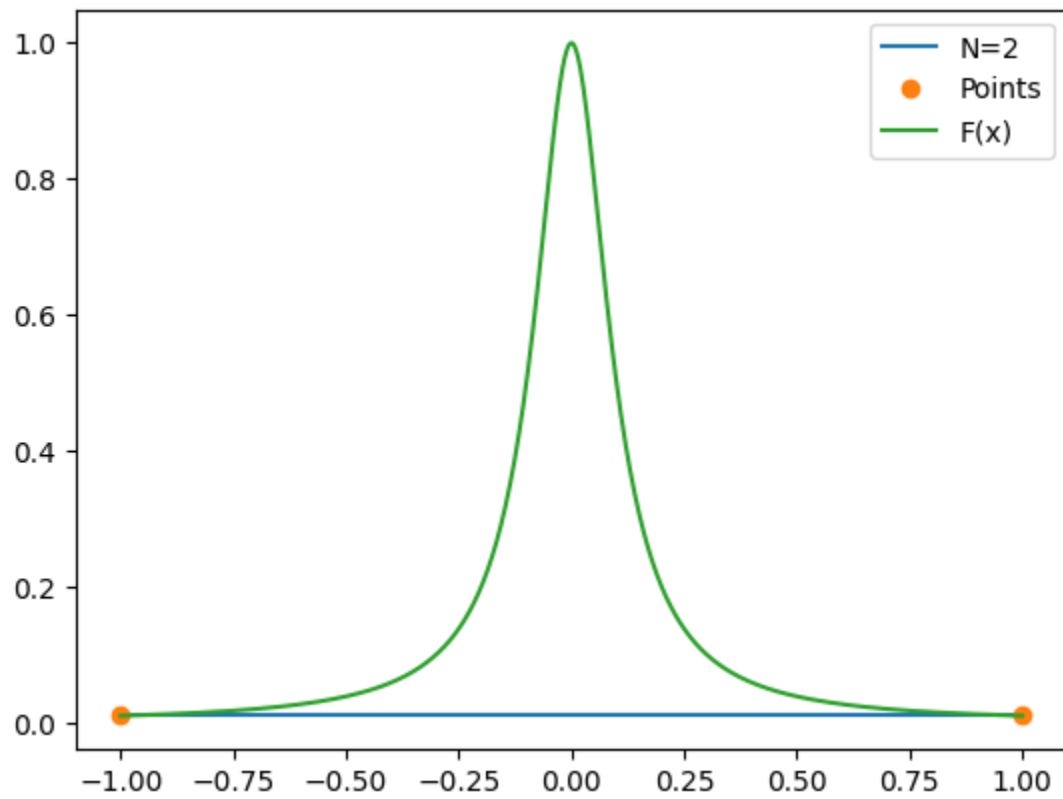
[[0.00990099]
 [0.          ]]

```

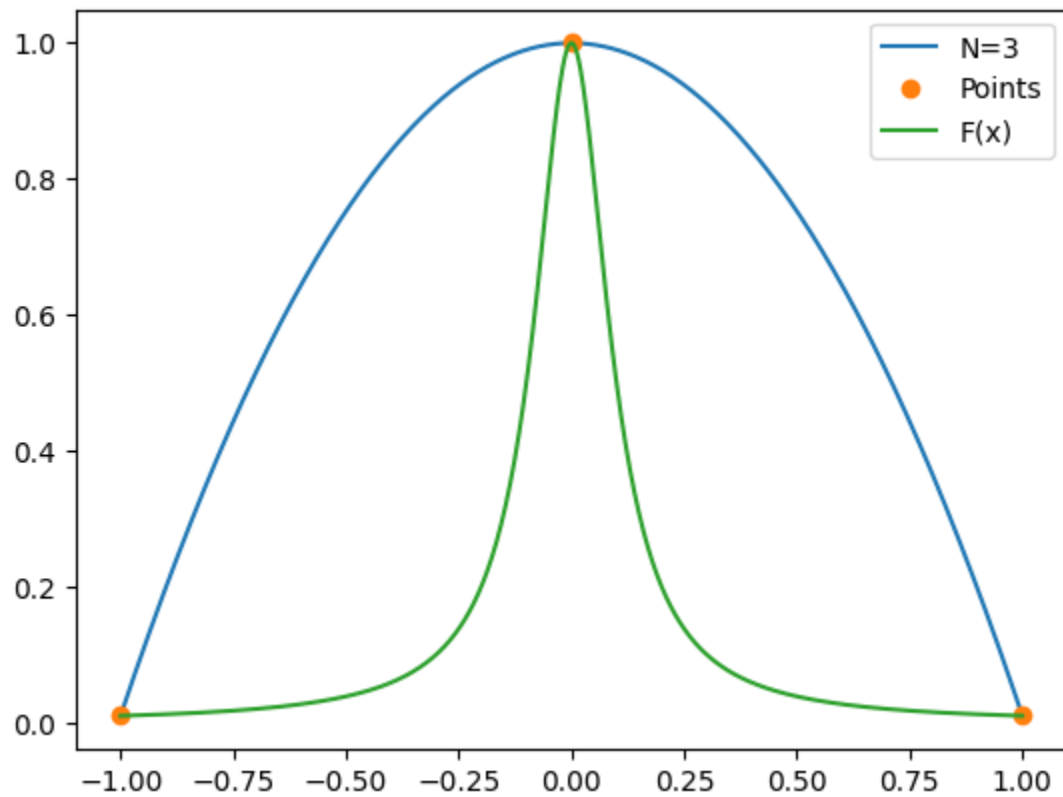
In [8]: `polyCoef = np.array(0)`
`poly=[0,0]`
`N2=2`
`xCont = np.linspace(-1,1,1001)`
`yCont = 1/(1+(10*xCont)**2)`
`while(max(poly)<100):`
`i2 = np.linspace(1,N2,N2)`
`x2= -1+(i2-1)*(2/(N2-1))`
`x2=x2.reshape((N2,1))`
`y2= 1/(1+(10*x2)**2)`

`polyCoef = vander(V1(x2,N2),y2)`
`poly = polyCoefficients(xCont,polyCoef)`
`print("At N=",N2,"The max value is:",max(poly))`
`plt.plot(xCont,poly, label = "N="+ str(N2))`
`plt.plot(x2,y2,"o", label = "Points")`
`plt.plot(xCont,yCont, label = "F(x)")`
`plt.legend()`
`plt.show()`
`plt.close()`
`N2+=1`

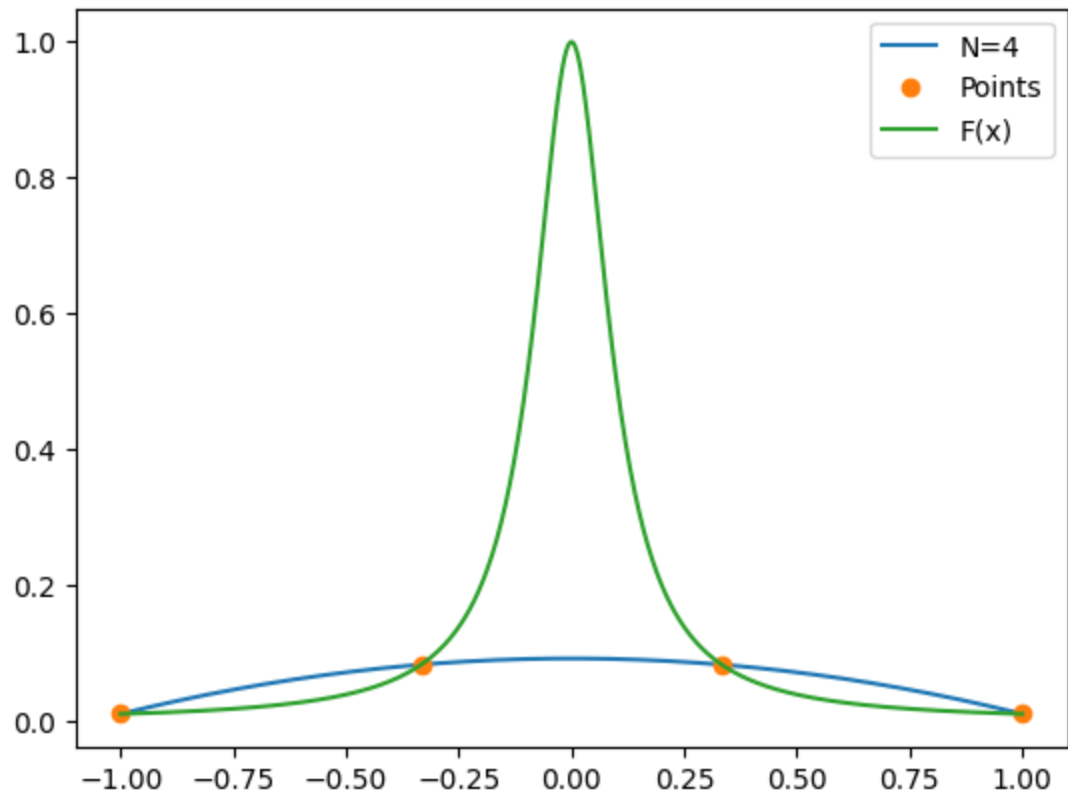
At $N = 2$ The max value is: 0.009900990099009901



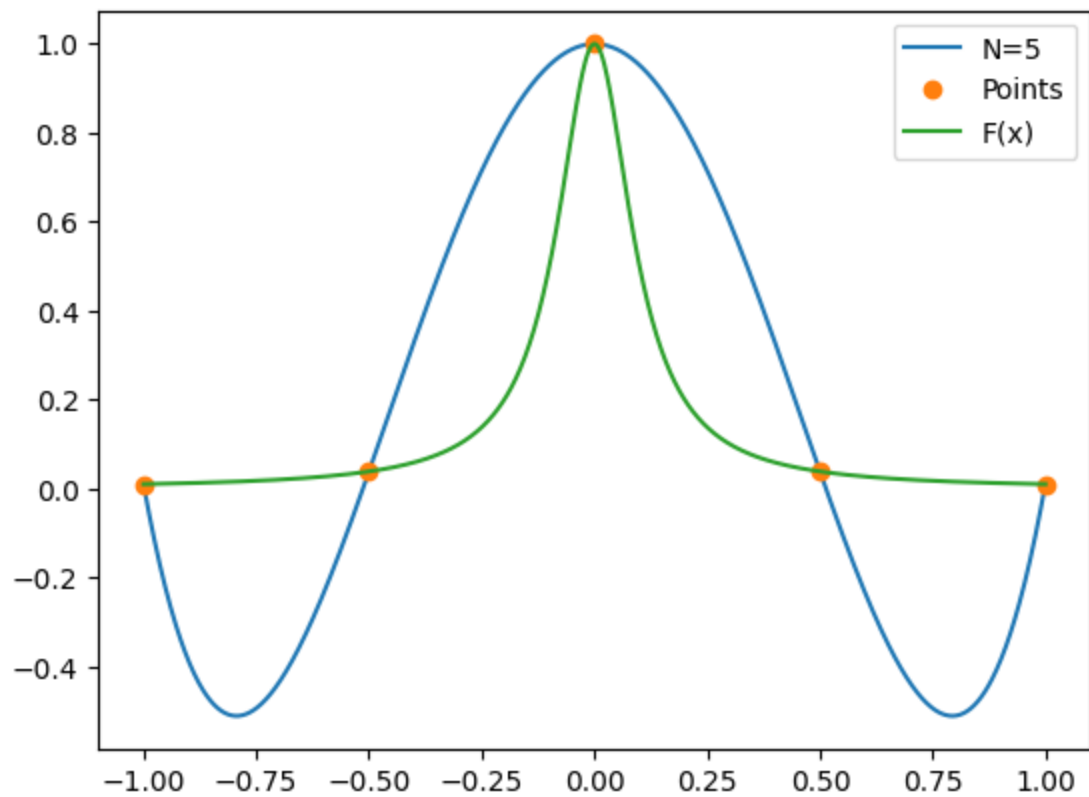
At $N = 3$ The max value is: 1.0



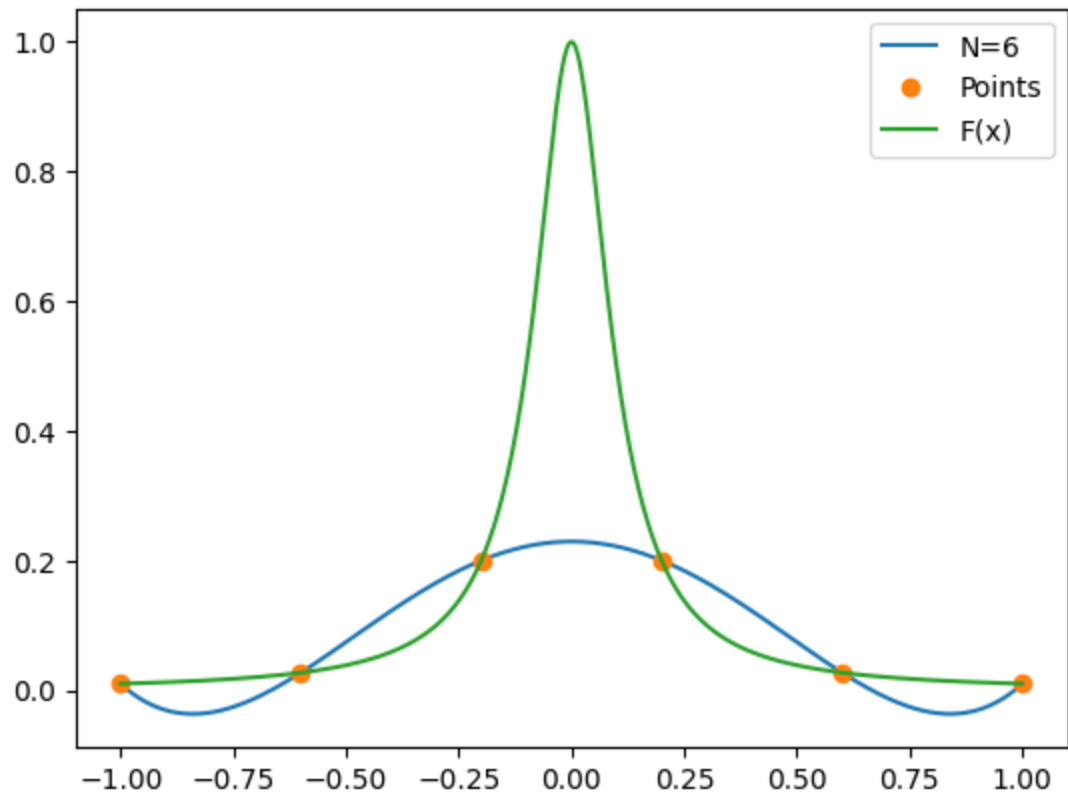
At $N = 4$ The max value is: 0.09165228449450451



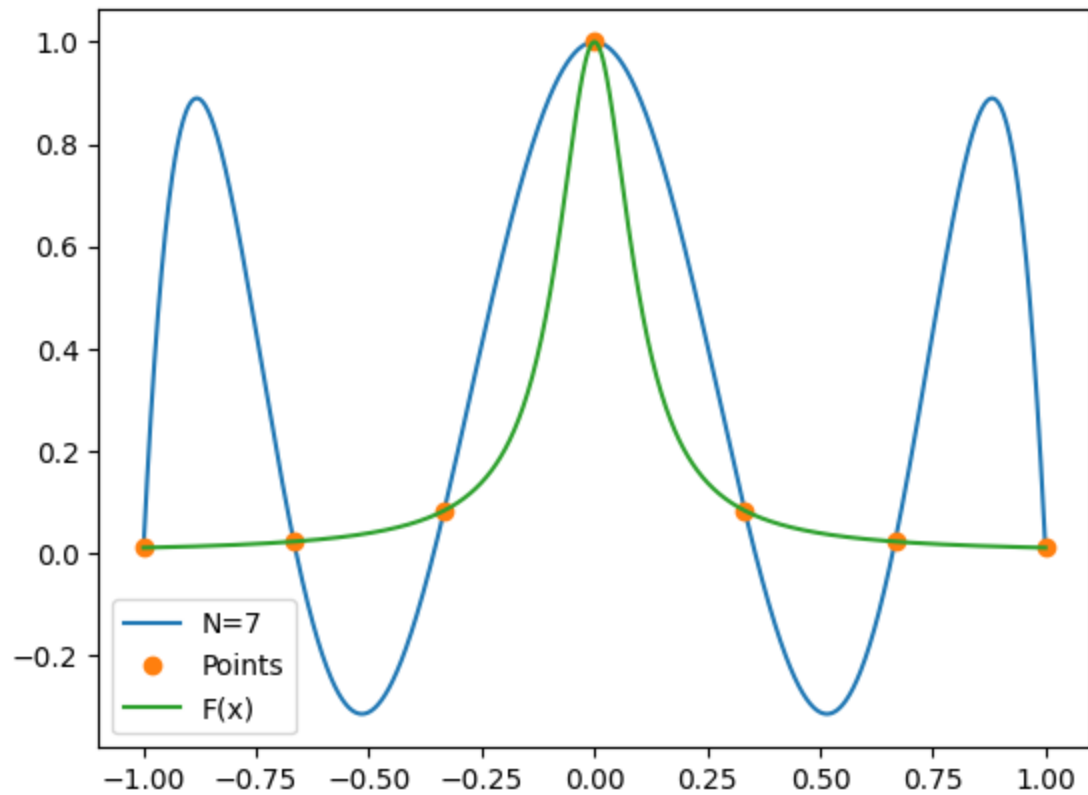
At $N=5$ The max value is: 1.0000000000000004



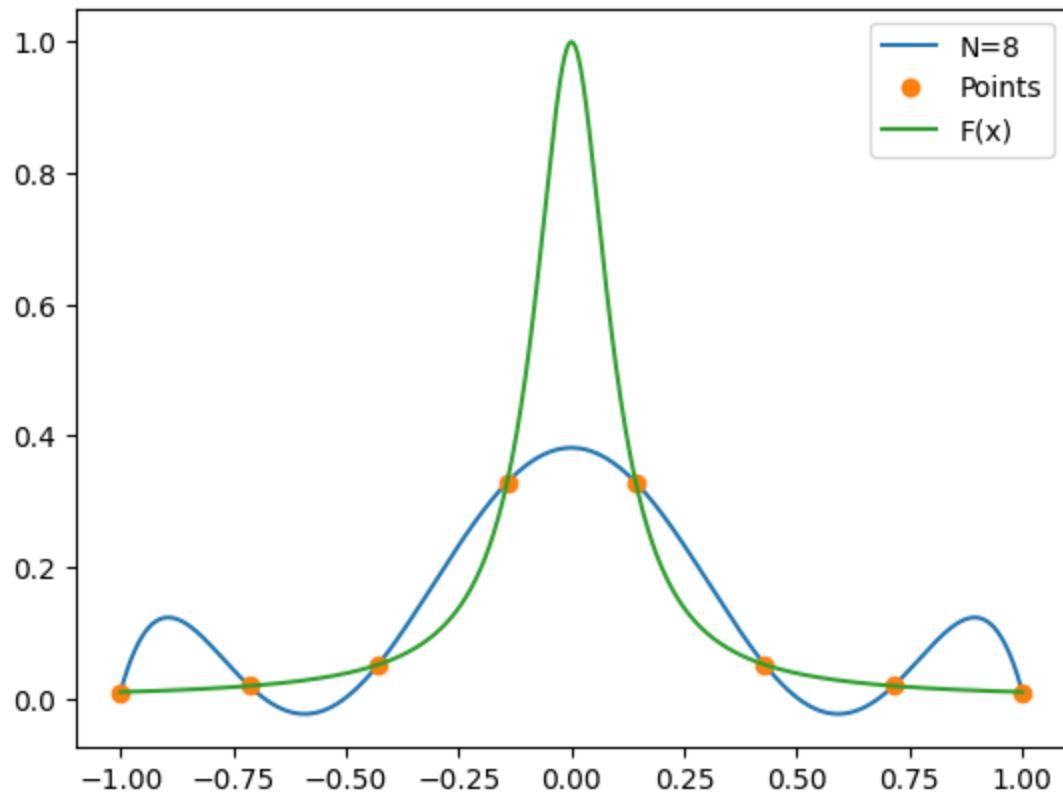
At $N=6$ The max value is: 0.22932833823922913



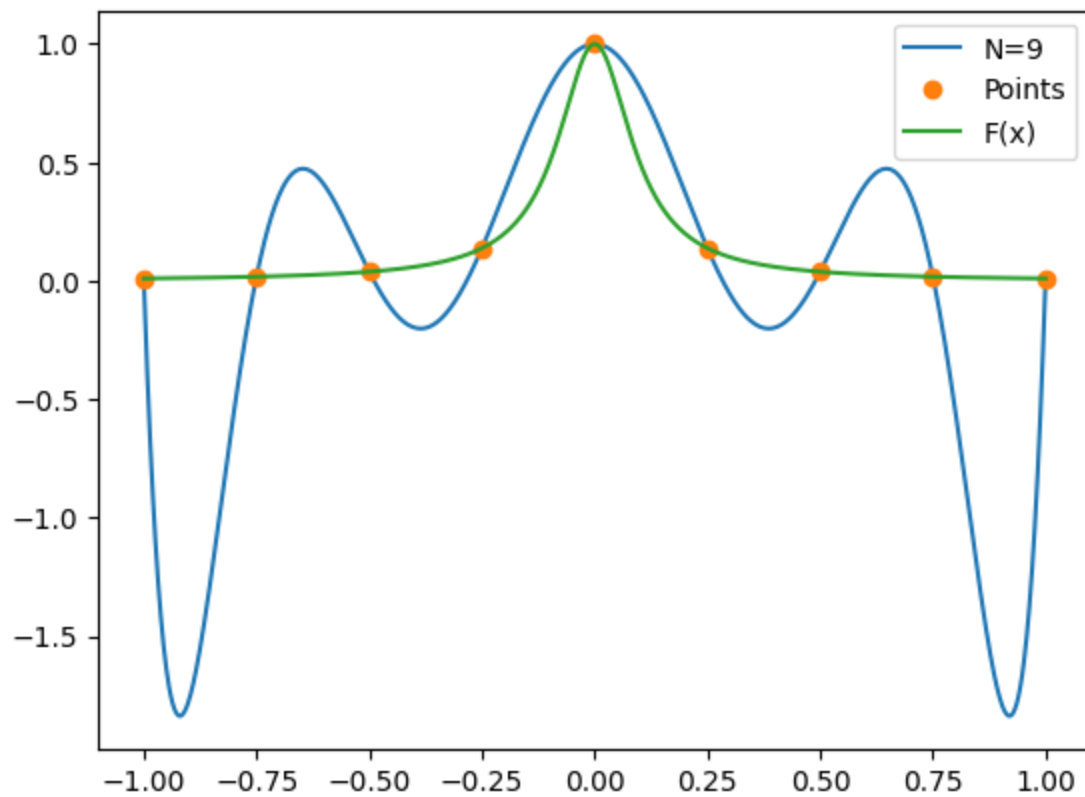
At $N=7$ The max value is: 0.9999999999999996



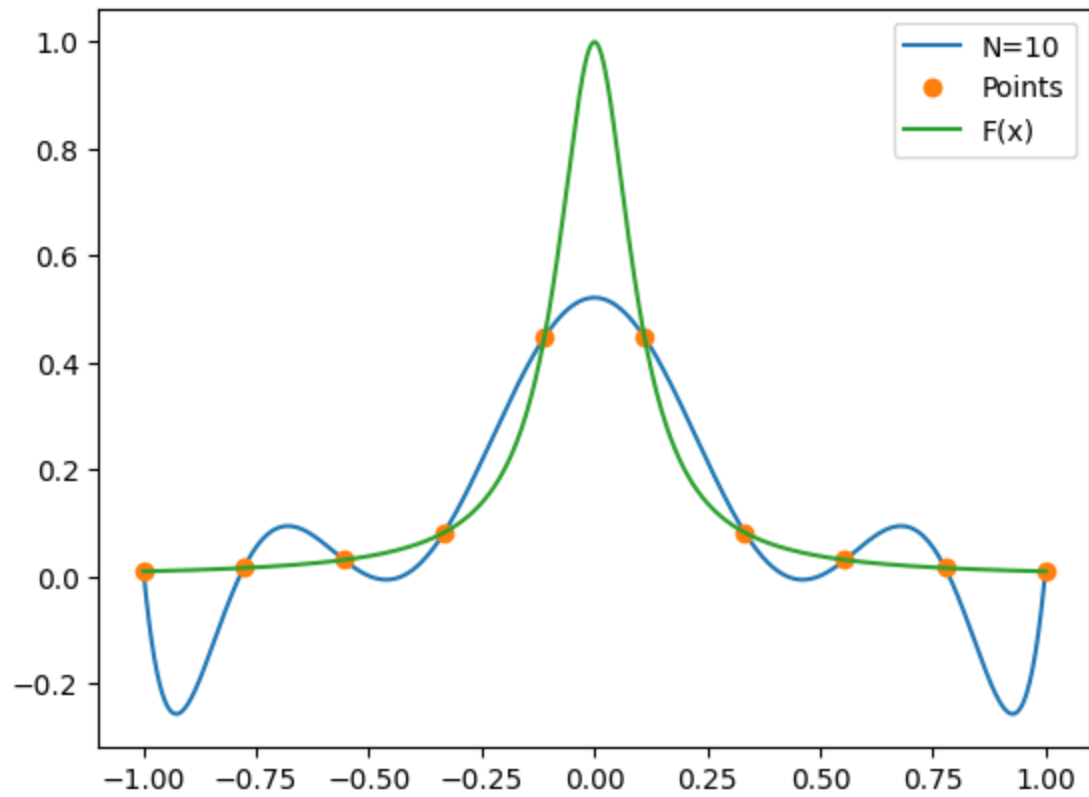
At $N=8$ The max value is: 0.3819283437466645



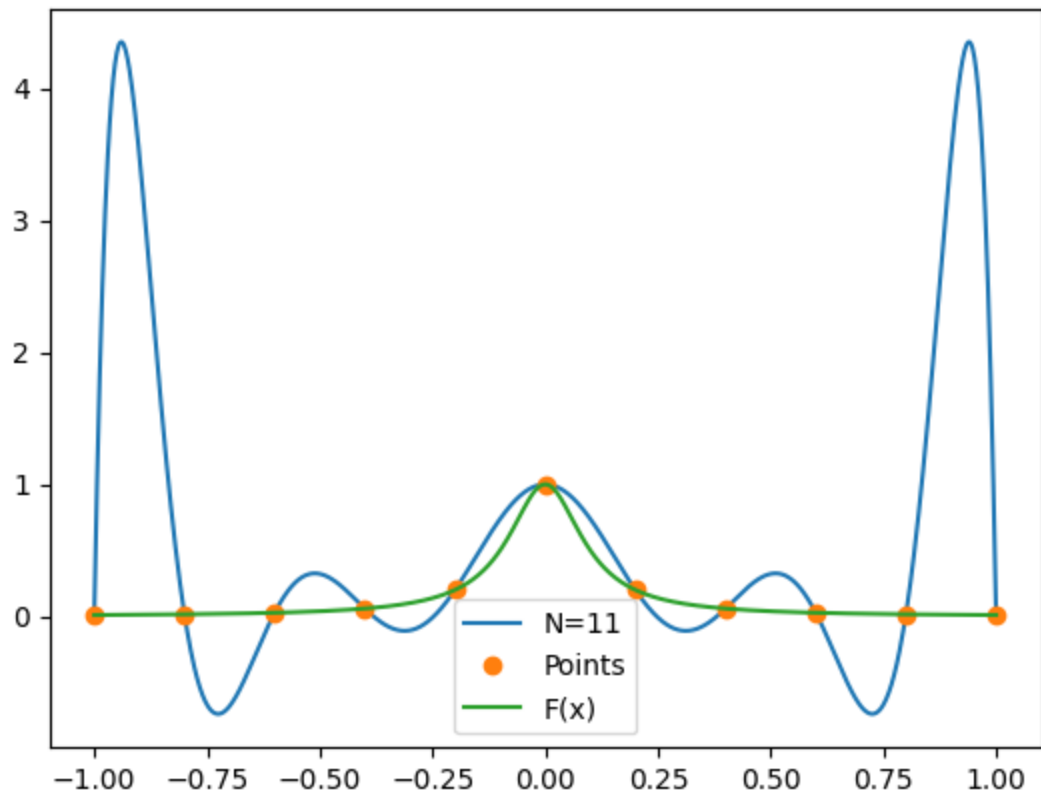
At $N=9$ The max value is: 0.9999999999999999



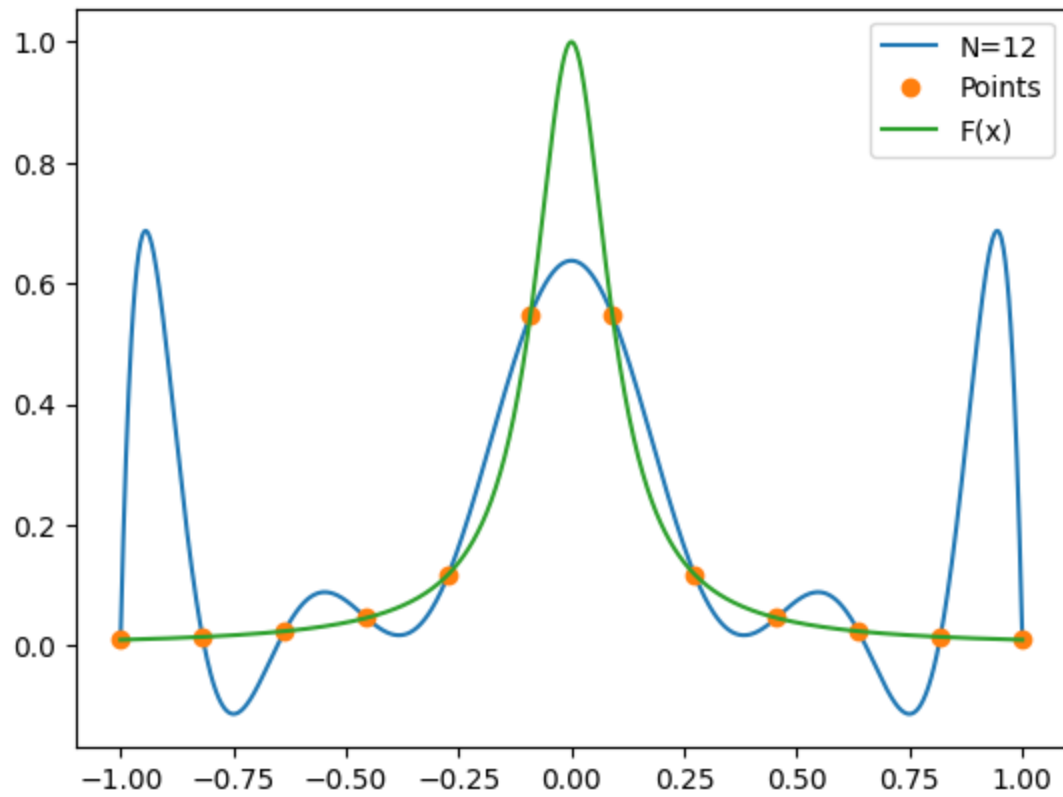
At $N=10$ The max value is: 0.5218049284237937



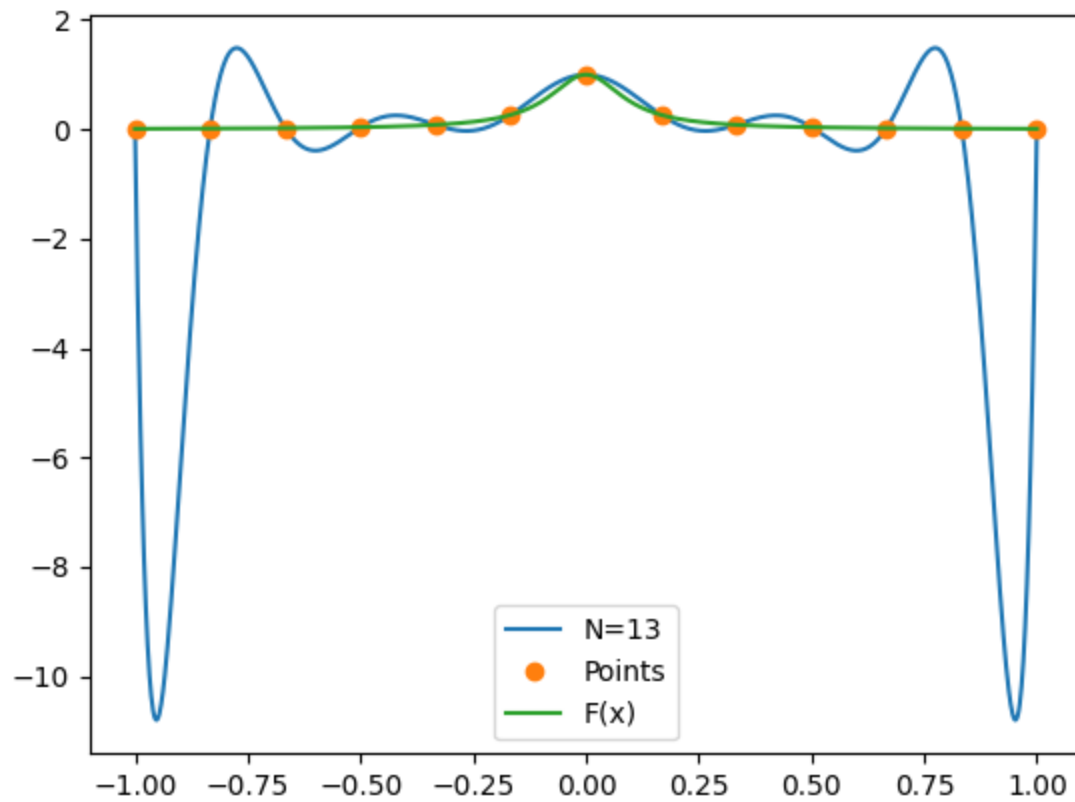
At $N=11$ The max value is: 4.351515169127424



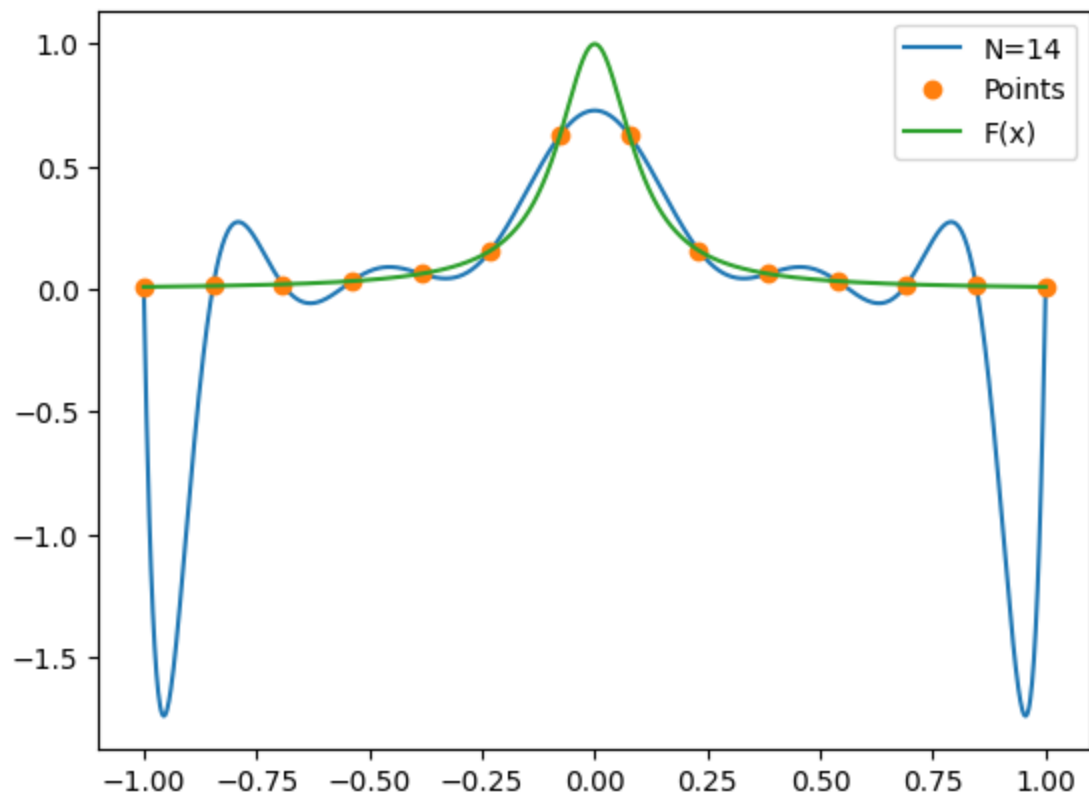
At $N=12$ The max value is: 0.6872844686330045



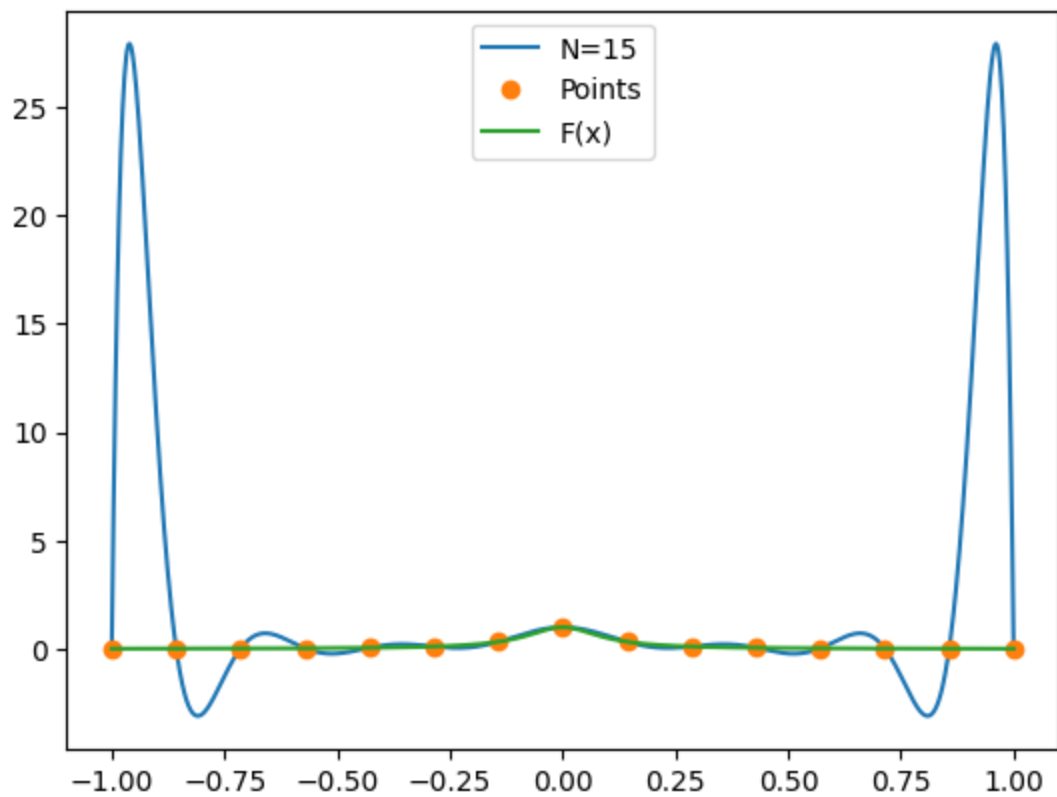
At $N=13$ The max value is: 1.4877279683334592



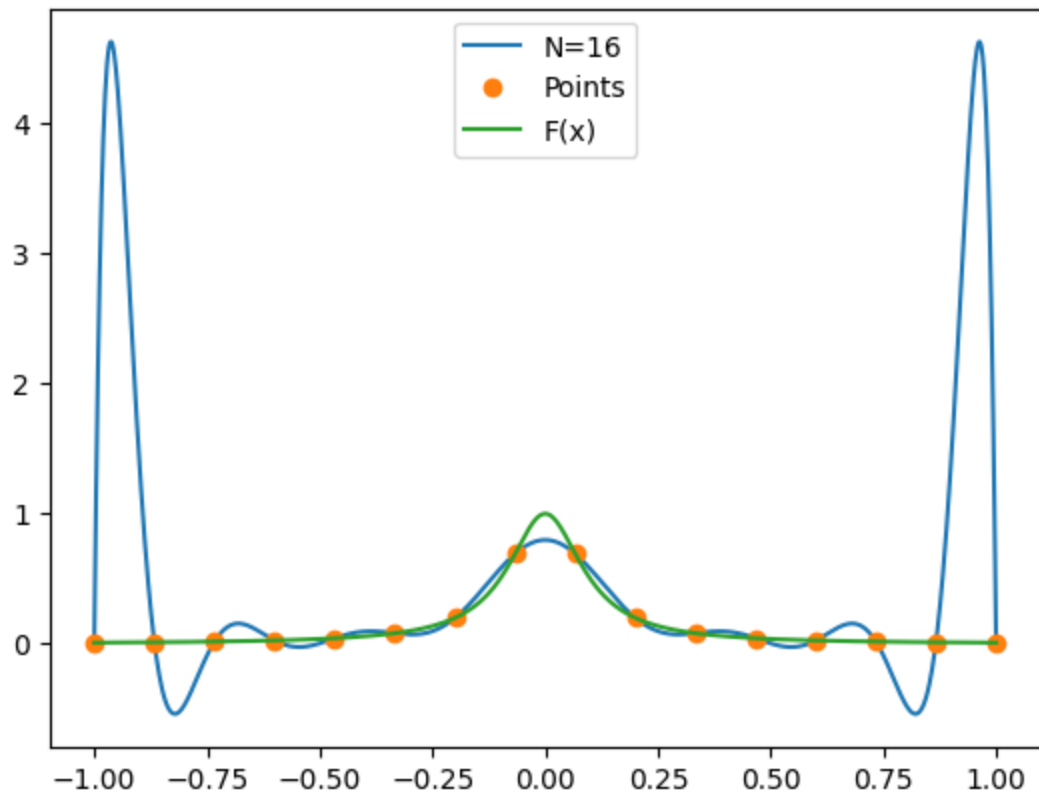
At $N=14$ The max value is: 0.7289405174432307



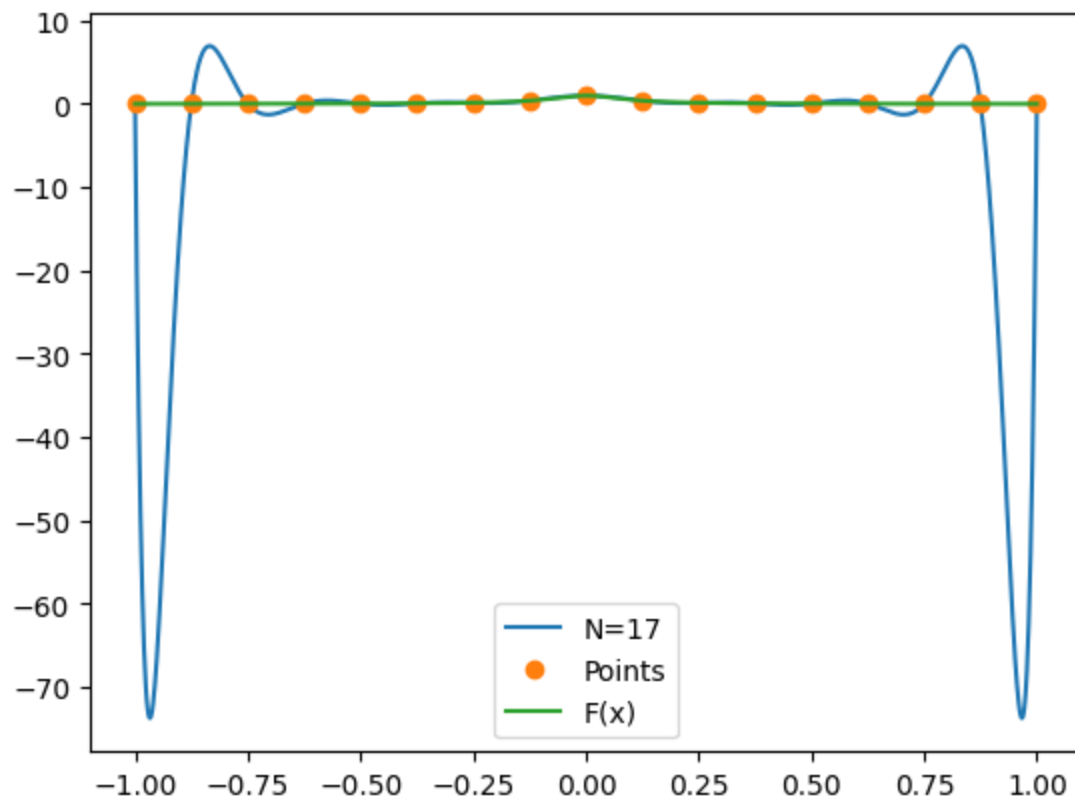
At $N = 15$ The max value is: 27.908207611999387



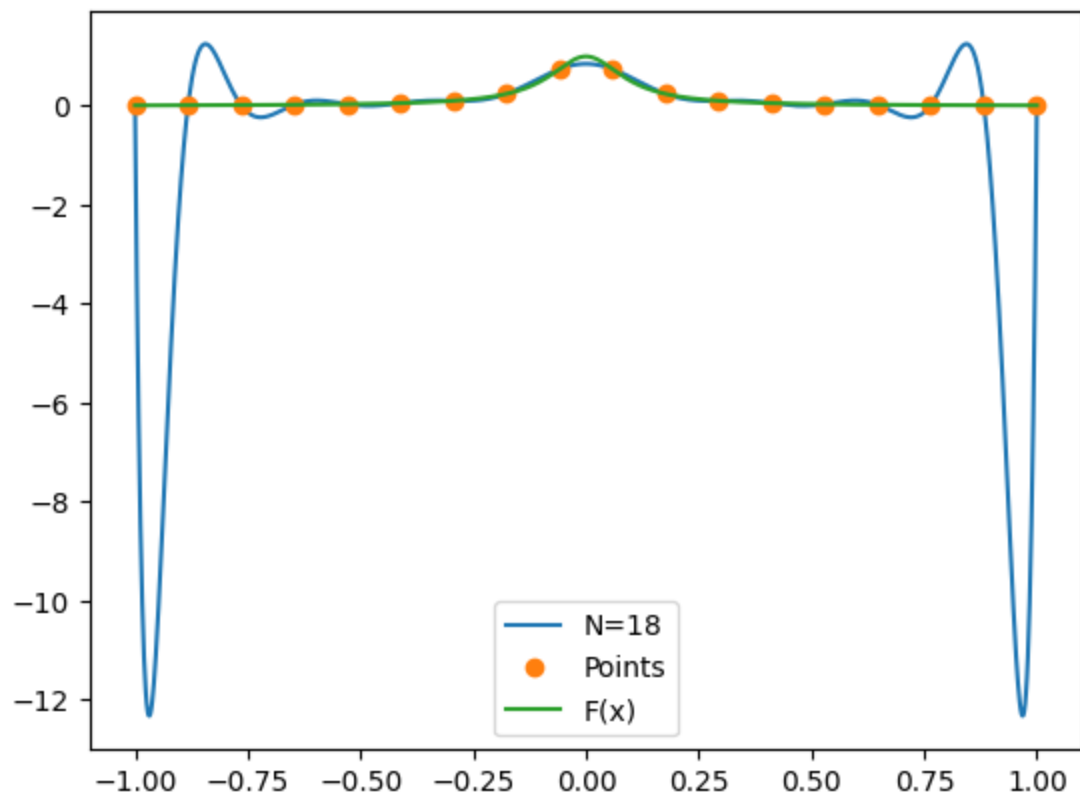
At $N = 16$ The max value is: 4.621874191009424



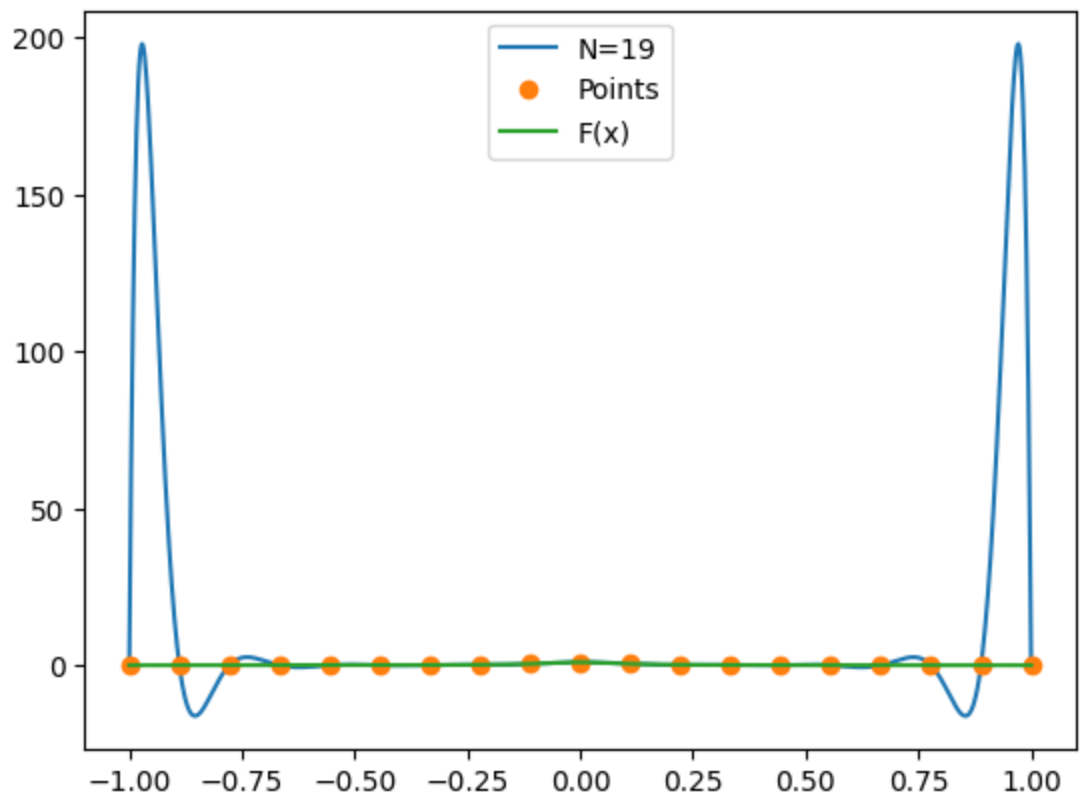
At $N = 17$ The max value is: 6.960564572897965



At $N = 18$ The max value is: 1.2514105362915586



At $N=19$ The max value is: 198.09925590845523



Here it is obvious that as N increases, the Runge Phenomenon increases in magnitude and creates issues with edge cases of the area. Eventually creating such large spikes that all interesting parts of the data have been lost as their magnitude is much less than the magnitude of variation at the edges.

Problem 2

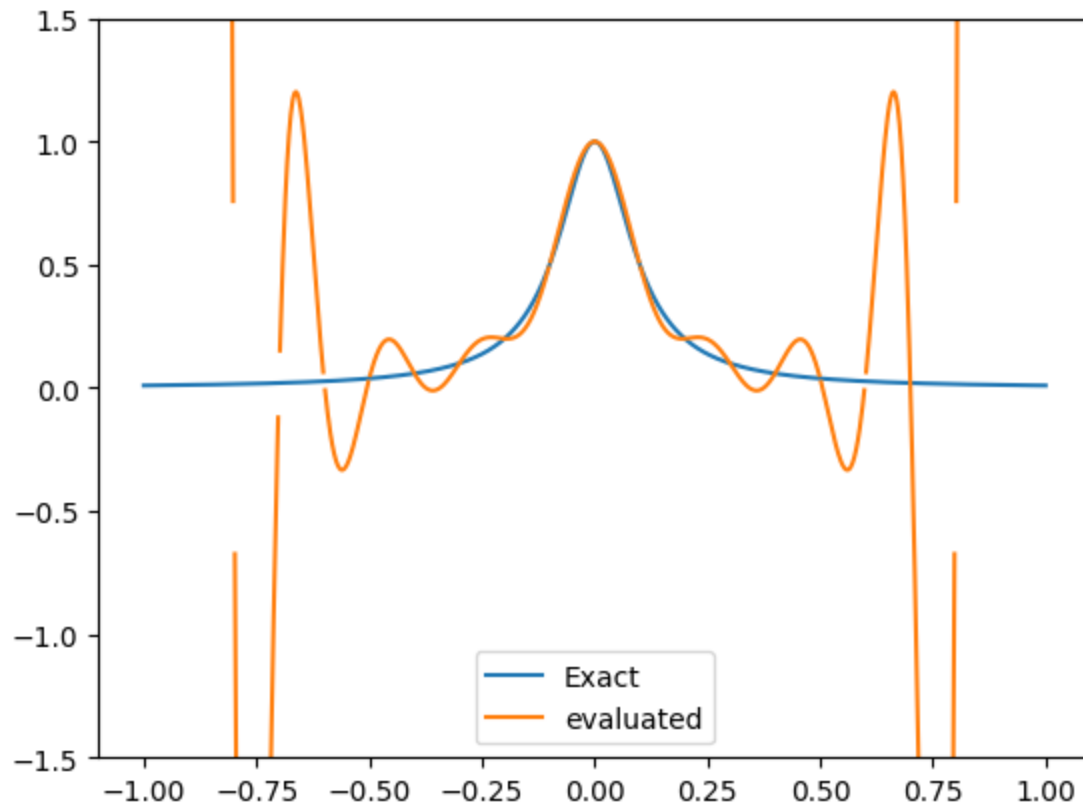
```
In [9]: """Evaluate Lagrange"""
def lagrangeEval(f,N,a,b):
    xint = np.linspace(a,b,N+1)
    yint = f(xint)

    Neval = 1000
    xeval = np.linspace(a,b,Neval+1)
    yeval_l= np.zeros(Neval+1)

    for i in range(Neval+1):
        yeval_l[i]= lag.eval_lagrange(xeval[i],xint,yint,N)
    y = f(xeval)
    plt.plot(xeval,y, label = "Exact")
    plt.plot(xeval,yeval_l, label = "evaluated")
    plt.ylim(-1.5,1.5)
    plt.legend()
```

```
In [10]: f = lambda x: 1/(1+(10*x)**2)
N=20
a=-1
b=1
lagrangeEval(f,N,a,b)
```

```
c:\Users\xman7\OneDrive - UCB-O365\Documents\APPM4600\Homework\Homework7\interp.py:2
4: RuntimeWarning: invalid value encountered in scalar divide
    peval = peval + phi*wj[j]*yint[j]/(xeval-xint[j])
```



This still exhibits the bad behaviour, but shows good behaviour for $x \ll |bound|$

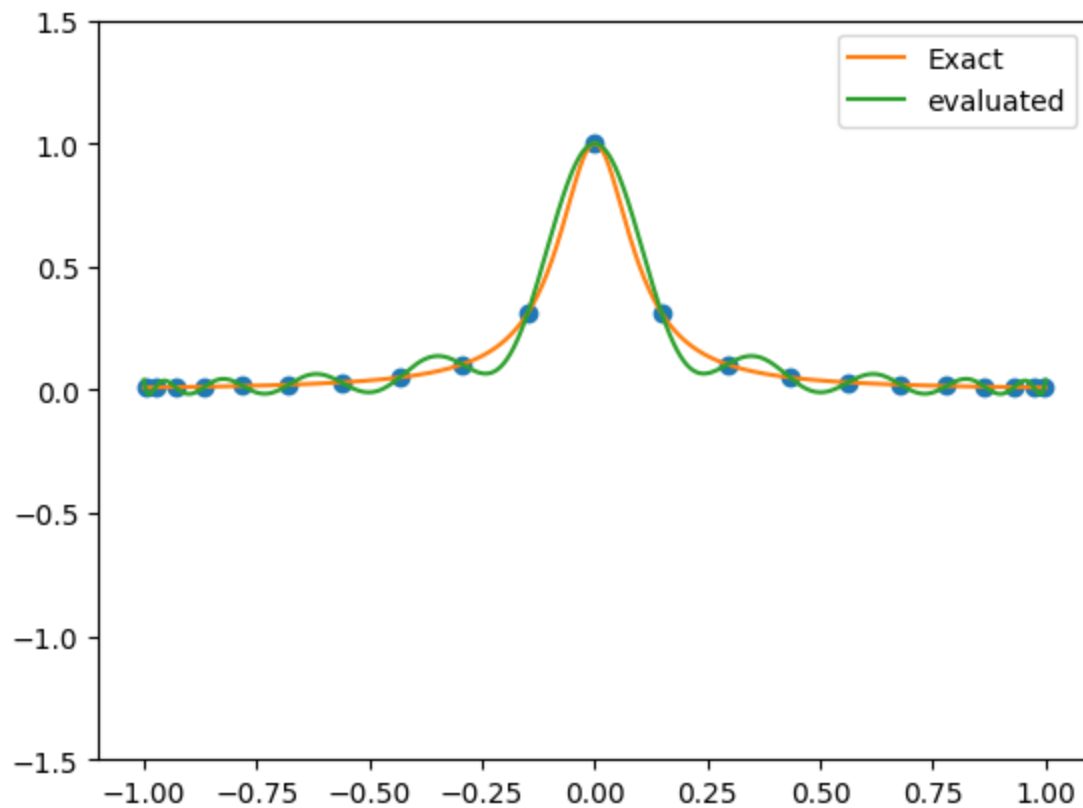
Problem 3

```
In [18]: def lagrangeEval2(f,N,a,b):
    jint = np.linspace(1,N,N+1)
    xint = np.cos(np.pi*(2*jint-1)/(2*N))
    yint = f(xint)
    plt.plot(xint,yint,'o')

    Neval = 1000
    xeval = np.linspace(a,b,Neval+1)
    yeval_l= np.zeros(Neval+1)

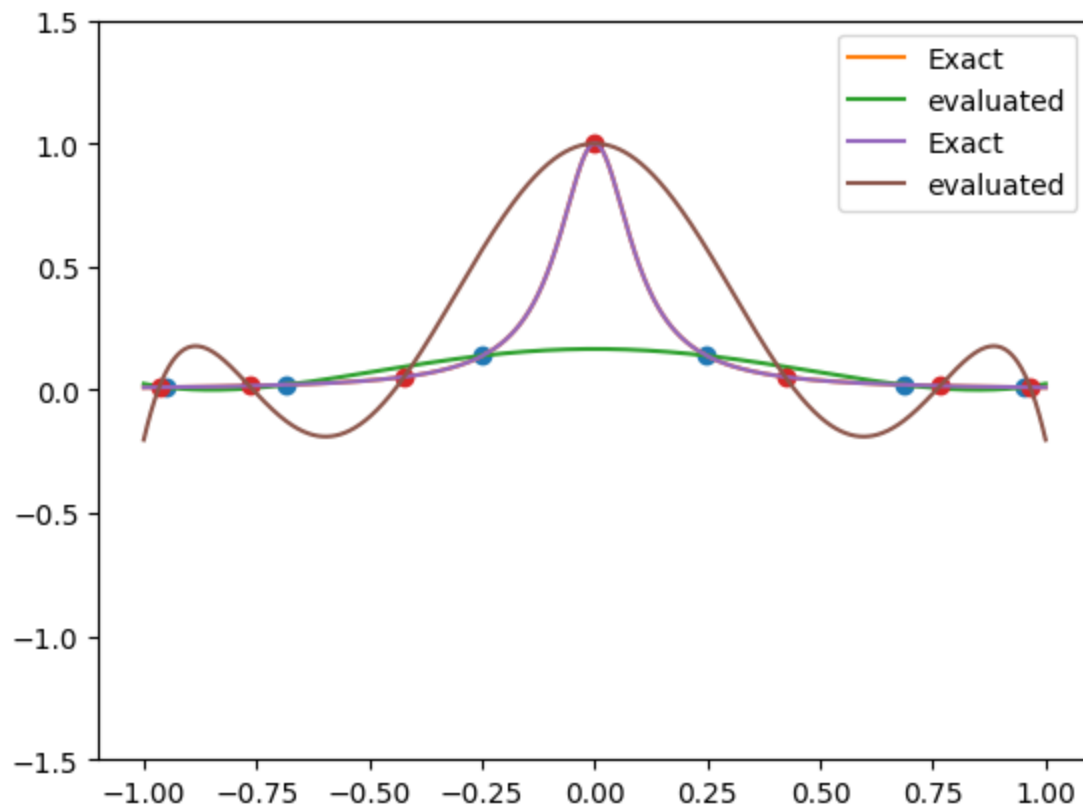
    for i in range(Neval+1):
        yeval_l[i]= lag.eval_lagrange(xeval[i],xint,yint,N)
    y = f(xeval)
    plt.plot(xeval,y, label = "Exact")
    plt.plot(xeval,yeval_l, label = "evaluated")
    plt.legend()
```

```
In [ ]: f = lambda x: 1/(1+(10*x)**2)
N=20
a=-1
b=1
lagrangeEval2(f,N,a,b)
```



Here you can see that using the chebyshev points the interpolation works wonderfully. So to get it to fail we will drastically decrease the number of points

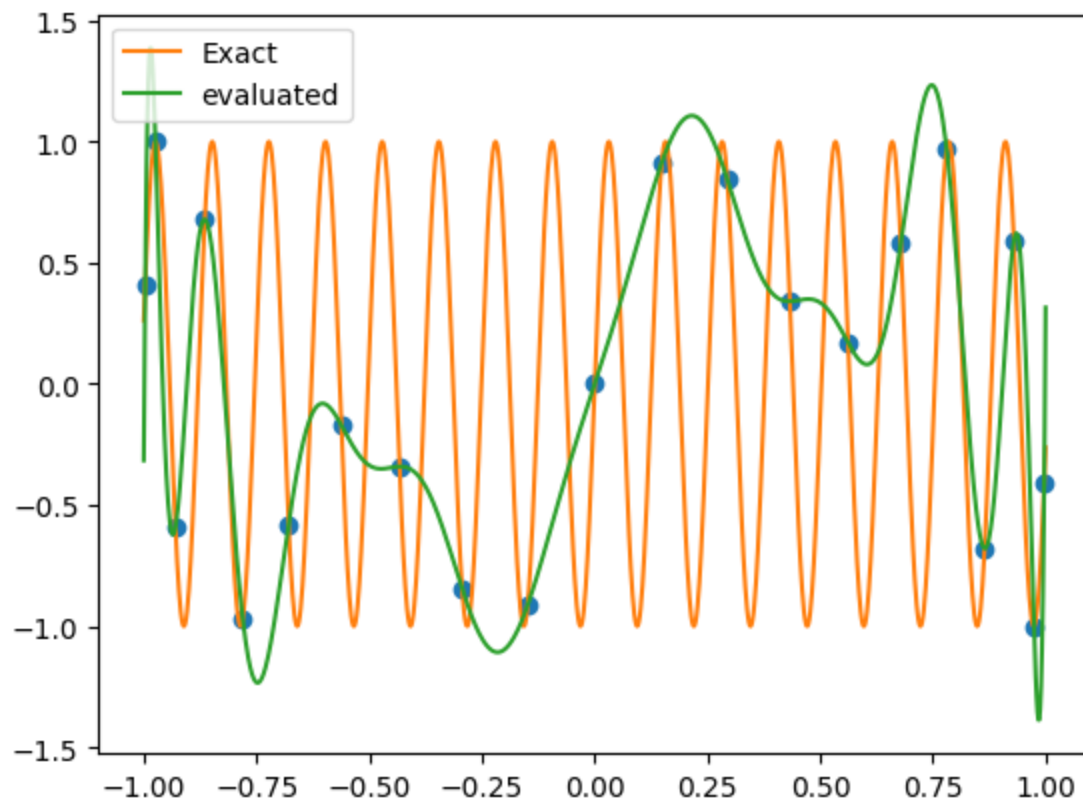
```
In [15]: f = lambda x: 1/(1+(10*x)**2)
N3=5
a=-1
b=1
lagrangeEval2(f,N3,a,b)
lagrangeEval2(f,N3+1,a,b)
```



This leads to a lack of accuracy for small x if the number of points is even, but for odd numbers it works fine.

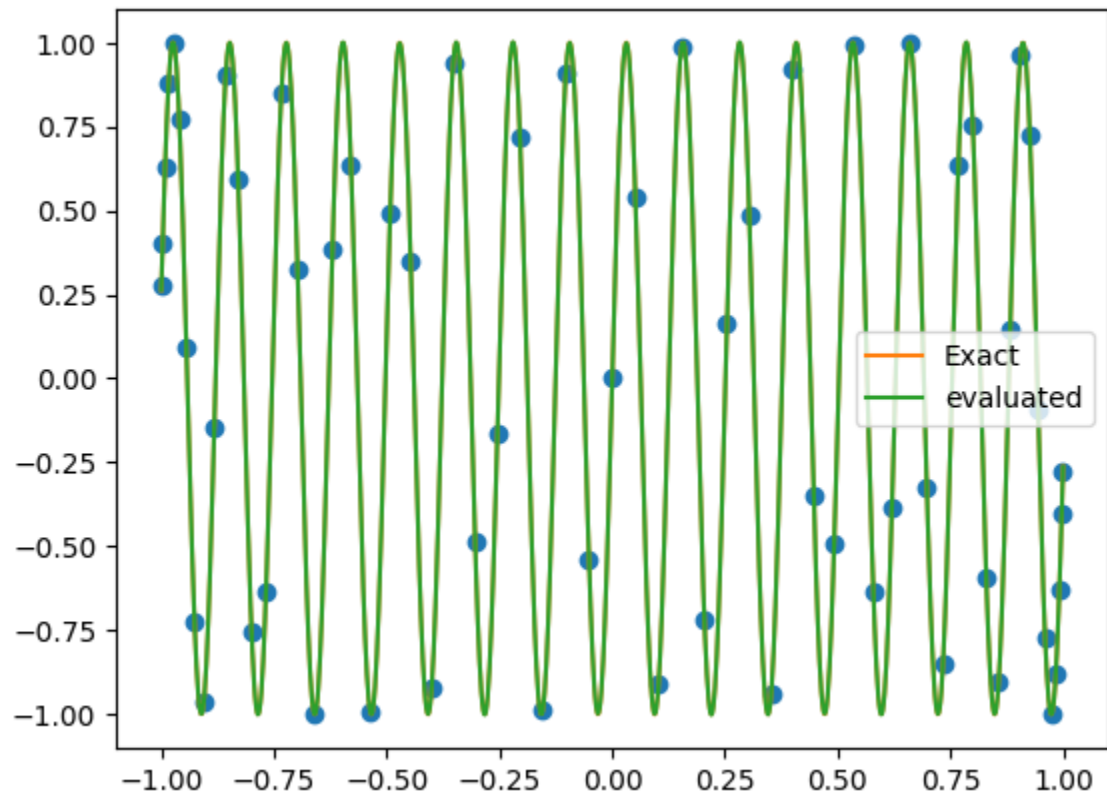
If we change f and its bounds, maybe the interpolation won't work as well:

```
In [27]: f2 = lambda x: np.sin(50*x)
lagrangeEval2(f2,N,a,b)
```



For a high frequency sinusoidal function this breaks for a reasonable N . You need $N \approx w$ where w is the frequency of your sinusoid.


```
In [29]: f2 = lambda x: np.sin(50*x)  
lagrangeEval2(f2,60,a,b)
```



```
In [ ]:
```