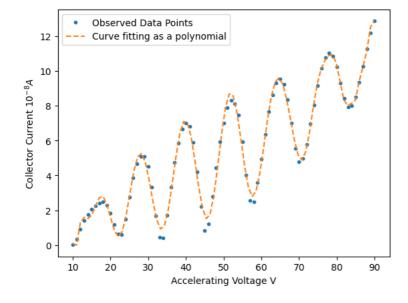
Franck-Hertz Experiment

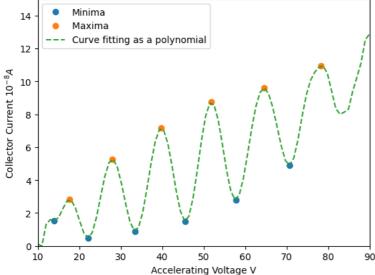
```
In [60]: %reset -f
          import numpy as np
import matplotlib.pyplot as plt
          from scipy.stats import linregress
          import pandas as pd
          import statistics
In [61]: df = pd.read_excel(r'/home/ashwin/Git/general-physics-lab/frank_hertz_2.xlsx')
              VG2K
                     CURRENT
          0
                        0.00
                  0
                        0.00
                  1
          1
2
3
                        0.00
                        0.00
                  3
          4
                  Δ
                        0.00
                        9.35
          86
                86
          87
                 87
                       10.25
          88
                 88
                       11.25
          89
                 89
                       12.18
          90
                 90
                       12.87
          [91 rows x 2 columns]
In [62]: VG2K=df['VG2K'].to_numpy()
          current=df['CURRENT'].to_numpy()
In [63]:
          x=VG2K[10:]
          y=current[10:]
          deg=80
          result=np.polyfit(x,y,deg)
          p = np.poly1d(result)
          xp = np.linspace(10, 90, 80)
          plt.figure()
          plt.plot(x, y,
                          '.',label="Observed Data Points")
          plt.plot(xp, p(xp),'--',label="Curve fitting as a polynomial")
plt.xlabel("Accelerating Voltage V")
          plt.ylabel("Collector Current $10^{-8}A$")
          plt.legend()
          /home/ashwin/.local/lib/python3.10/site-packages/numpy/lib/polynomial.py:666: RuntimeWarning: overflow encountered in multipl
```

scale = NX.sqrt((lhs*lhs).sum(axis=0))
/home/ashwin/.local/lib/python3.10/site-packages/IPython/core/interactiveshell.py:3378: RankWarning: Polyfit may be poorly conditioned
 exec(code_obj, self.user_global_ns, self.user_ns)

Out[63]: <matplotlib.legend.Legend at 0x7f8e5c78c280>



```
In [64]: crit = p.deriv().r
            r_crit = crit[crit.imag==0].real
            test = p.deriv(2)(r\_crit)
            # compute local minima
            # excluding range boundaries
            x_min = r_crit
            x_min=x_min[::-1]
            x_min=sorted(x_min)
            #Removing the ourlier values.
            x_{\min}=x_{\min}[2:-1]
            y_min = p(x_min)
            plt.plot( x_min[::2], y_min[::2], 'o' , label="Minima" ) plt.plot( x_min[1::2], y_min[1::2], 'o' , label="Maxima " )
            plt.plot(xp, p(xp),'--',label="Curve fitting as a polynomial")
plt.xlabel("Accelerating Voltage V")
plt.ylabel("Collector Current $10^{-8}A$")
            plt.xlim([10,90])
            plt.ylim([0,15])
            plt.legend()
            plt.show()
```



```
In [65]: diff=[]
        for i in range(1,len(x_min)-2):
          diff.append(abs(x_min[i]-x_min[i+2]))
        print(diff)
       print(len(diff))
        408674048408, 12.925463151598443, 13.769943646212766]
In [66]: statistics.stdev(diff)
Out[66]: 0.9726228575481425
In [67]: statistics.mean(diff)
Out[67]: 12.130329522412062
In [68]: print("Percentage error compared with literature is \n {0}%".format(round(abs((statistics.mean(diff)-11.83)/11.83*100),2)))
        Percentage error compared with literature is
        2.54%
In [69]: print("The Energy drop on inelastic collision of electron on Argon atom is : : \n {0} +- {1} ".format(round(statistics.mean(
       The Energy drop on inelastic collision of electron on Argon atom is : : 12.13 +- 0.97
```

This Concludes and quantization of energy in atomic model as per Quantum theory.