

ECSE 543: Numerical Methods
Assignment 3



Student Name: Dafne Culha
Student ID: 260785524

Department of Electrical and Computer Engineering
McGill University, Canada
November, 2020

1. You are given a list of measured BH points for M19 steel (Table 1), with which to construct a continuous graph of B versus H.

B (T)	H (A/m)
0.0	0.0
0.2	14.7
0.4	36.5
0.6	71.7
0.8	121.4
1.0	197.4
1.1	256.2
1.2	348.7
1.3	540.6
1.4	1062.8
1.5	2318.0
1.6	4781.9
1.7	8687.4
1.8	13924.3
1.9	22650.2

Table 1: B-H Data for M19 Steel

(a) Interpolate the first 6 points using full-domain Lagrange polynomials. Is the result plausible, i.e. do you think it lies close to the true B versus H graph over this range?

To interpolate specified points, algorithm discussed in class was followed. To interpolate 6 points, 6 Lagrange polynomials of 5th order are needed. Each of these polynomials are given by

$$L_j(x) = \frac{F_j(x)}{F_j(x_j)}$$

and where

$$F_j(x) = \prod_{r=1, r \neq j}^n (x - x_r)$$

Finally, using these, $y(x)$ can be approximated as

$$y(x) = \sum_{j=1}^{n=6} a_j \cdot L_j$$

where $a_j = y(x_j)$.

The whole implementation can be seen in interpolation.py. By using this program, the first 6 points were interpolated using full-domain Lagrange polynomials. The interpolated function was found to be $H = 14.062500000009 \cdot B^{**5} - 963.541666666669 \cdot B^{**4} + 873.4375000000015 \cdot B^{**3} - 215.208333333338 \cdot B^{**2} + 88.65000000000005 \cdot B$ as it is displayed in Figure 1.1.1. As expected, the function is a 5th degree polynomial. This function was plotted, and it's displayed in Figure 1.1.2. As it can be seen, this graph resembles a true B vs. H graph. Thus, it's concluded that the result is plausible.

```
-- /Users/dafneculha/Desktop/A3-260785524/interpolation.py
```

Part A

```
B = [0.0, 0.2, 0.4, 0.6, 0.8, 1.0]
```

```
H = [0.0, 14.7, 36.5, 71.7, 121.4, 197.4]
```

```
H = 414.062500000009*B**5 - 963.54166666669*B**4 + 873.437500000015*B**3 - 215.208333333338*B**2 + 88.6500000000005*B
```

Figure 1.1.1: Interpolation of First 6 Points

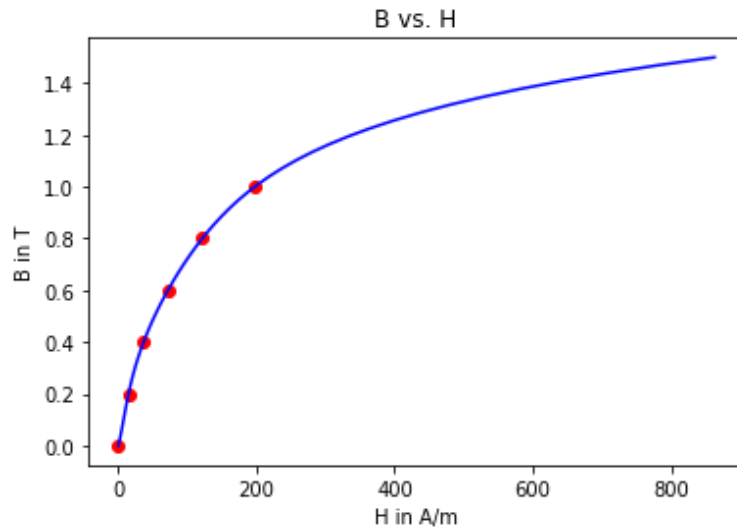


Figure 1.1.2: B vs. H

(b) Now use the same type of interpolation for the 6 points at $B = 0, 1.3, 1.4, 1.7, 1.8, 1.9$. Is this result plausible?

The data at the specified points was interpolated using the program written for previous part again. The interpolated function was found to be $H = 156393.280524088*B^{**5} - 966235.57224511*B^{**4} + 2253820.22115058*B^{**3} - 2337828.82945774*B^{**2} + 906781.854422079*B$ as it is displayed in Figure 1.2.1. This function was plotted, and it's displayed in Figure 1.2.2. As it can be seen, this graph has too many wiggles and does not resemble the B vs. H graph. Thus, it's concluded that the result is not plausible.

Part B

```
B = [0.0, 1.3, 1.4, 1.7, 1.8, 1.9]
```

```
H = [0.0, 540.6, 1062.8, 8687.4, 13924.3, 22650.2]
```

```
H = 156393.280524088*B**5 - 966235.57224511*B**4 + 2253820.22115058*B**3 - 2337828.82945774*B**2 + 906781.854422079*B
```

Figure 1.2.1: Interpolation of Data at $B = 0, 1.3, 1.4, 1.7, 1.8, 1.9$ with full-domain Lagrange polynomials

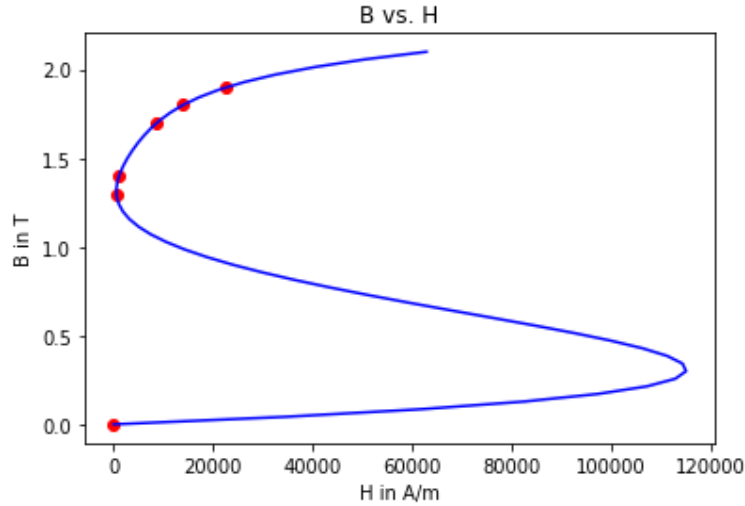


Figure 1.2.2: B vs. H

(c) An alternative to full-domain Lagrange polynomials is to interpolate using cubic Hermite polynomials in each of the 5 subdomains between the 6 points given in (b). With this approach, there remain 6 degrees of freedom - the slopes at the 6 points. Suggest ways of fixing the 6 slopes to get a good interpolation of the points. Test your suggestion and comment on the results.

$$L_j(x) = \frac{F_j(x)}{F_j(x_j)}$$

$$L_j'(x) = \frac{dL_j}{dx}$$

$$U_j(x) = [-1 - 2L_j'(x_j)(x - x_j)] \cdot L_j(x) \cdot L_j(x)$$

$$V_j(x) = (x - x_j) \cdot L_j(x) \cdot L_j(x)$$

where $U_j(x)$ and $V_j(x)$ are polynomials of degree $2n-1 = 11$.

Finally, using these, $y(x)$ can be obtained as

$$y(x) = \sum_{j=1}^{n=6} a_j \cdot U_j(x) + b_j \cdot V_j(x)$$

where $a_j = y(x_j)$

and

$$b_j = y'(x_j) = \frac{y(x_{j+1}) - y(x_j)}{x_{j+1} - x_j}$$

The whole implementation can be seen in interpolation.py. Using this program, the data at the specified points was interpolated using cubic Hermite polynomials. The interpolated function was found to be $H =$

$1734143651.0417*B^{11} - 25207926894.739*B^{10} + 162399433091.79*B^9 - 608575443825.474*B^8 + 1461887595692.59*B^7 - 2334363871169.64*B^6 + 2477827583874.83*B^5 - 1685862719412.73*B^4 + 667146472644.371*B^3 - 116995129462.297*B^2 + 415.846153846154*B$ as it is presented in Figure 1.3.1. As expected, the function is a 11th degree polynomial. This function was plotted, and it's displayed in Figure 1.3.2. As it can be seen, this graph resembles the B vs. H graph. Thus, it's concluded that the result is plausible.

Part C

$B = [0.0, 1.3, 1.4, 1.7, 1.8, 1.9]$
 $H = [0.0, 540.6, 1062.8, 8687.4, 13924.3, 22650.2]$
 $H = 1734143651.0417*B^{11} - 25207926894.739*B^{10} + 162399433091.79*B^9 - 608575443825.474*B^8 + 1461887595692.59*B^7 - 2334363871169.64*B^6 + 2477827583874.83*B^5 - 1685862719412.73*B^4 + 667146472644.371*B^3 - 116995129462.297*B^2 + 415.846153846154*B$

Figure 1.3.1: Interpolation of Data at $B = 0, 1.3, 1.4, 1.7, 1.8, 1.9$ with Cubic Hermite polynomials

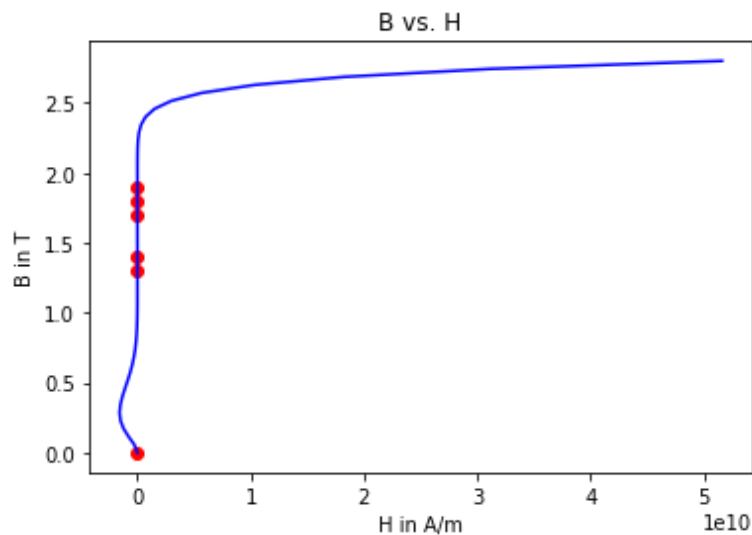


Figure 1.3.2: B vs. H

(d) The magnetic circuit of Figure 1 has a core made of M19 steel, with a cross-sectional area of 1 cm^2 . $L_C=30\text{cm}$ and $L_A=0.5\text{cm}$. The coil has $N=800$ turns and carries a current $I=10\text{A}$. Derive a (nonlinear) equation for the flux ψ in the core, of the form $f(\psi) = 0$.

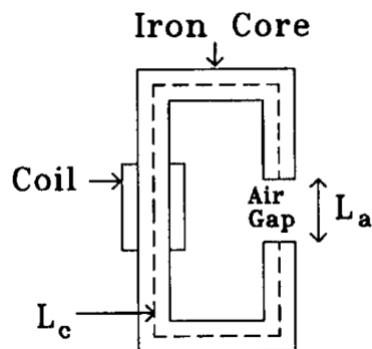


Figure 1.4.1: M19 Steel

By using knowledge learnt in Power Engineering and revised in ECSE 543 lectures,

$$MMF = \mathcal{F} = N \cdot I = R_{Total} \times \psi$$

$$R_{Total} = R_A + R_C$$

$$R_A = \frac{L_A}{\mu_A \cdot A_A}$$

$$\mu_c = \frac{B_C}{H_C}$$

$$B_c = \frac{\psi}{A_c}$$

$$R_c = \frac{L_c}{\mu_c \cdot A_c} = \frac{H_C}{B_C} \cdot \frac{L_c}{A_c} = \frac{H_C}{\psi} \cdot L_c$$

$$\mathcal{F} = N \cdot I = 800 \text{ t} \cdot 10A = 8000 \text{ A} \cdot \text{t}$$

$$\mu_A = \mu_0 = 4\pi \times 10^{-7}$$

$$R_A = \frac{L_A}{\mu_A \cdot A_A} = \frac{0.5 \times 10^{-2}}{4\pi \times 10^{-7} \times 1 \times 10^{-4}} = 39\,788\,735.77$$

$$R_c = \frac{L_c}{\mu_c \cdot A_c} = \frac{H_C(\psi)}{B_C} \cdot \frac{L_c}{A_c} = \frac{H_C(\psi)}{\psi} \cdot L_c = \frac{H_C(\psi)}{\psi} \cdot 30 \times 10^{-2} = 0.3 \times \frac{H_C(\psi)}{\psi}$$

$$R_{Total} = R_A + R_c = \frac{H_C(\psi)}{\psi} \cdot 30 \times 10^{-2} + 39\,788\,735.77$$

$$\mathcal{F} = N \cdot I = R_{Total} \times \psi$$

$$8000 = \psi \cdot \left(\frac{H_C(\psi)}{\psi} \cdot L_c + R_A \right) = H_C \times 0.3 + 39\,788\,735.77 \cdot \psi$$

$$f(\psi) = 0 = H_C(\psi) \cdot L_c + R_A \cdot \psi - N \cdot I$$

$$f(\psi) = 0 = H_C(\psi) \times 0.3 + 39\,788\,735.77 \cdot \psi - 8000$$

(e) Solve the nonlinear equation using Newton-Raphson. Use a piecewise-linear interpolation of the data in Table 1. Start with zero flux and finish when $|\hat{f}(\psi) / \hat{f}(0)| < 10^{-6}$. Record the final flux, and the number of steps taken.

To use Newton-Raphson method, first $f(\psi)$ and $f'(\psi)$ need to be found.

$$f(\psi) = 0 = H_C(\psi) \times 0.3 + 39\,788\,735.77 \cdot \psi - 8000 = H_C(\psi) \cdot L_C + R_A \cdot \psi - 8000$$

and

$$f'(\psi) = 0 = H'_C(\psi) \times 0.3 + 39\,788\,735.77 = H'_C(\psi) \cdot L_C + R_A$$

To get $H_C(\psi)$ and $H'_C(\psi)$ piecewise linear approximation of data in Table 1 was used. To find these,

$$\mu_c = \frac{\Delta B_c}{\Delta H_c} = \frac{B_i - B_{i-1}}{H_i - H_{i-1}}$$

$$B_c(\psi) = \psi / \text{Area}$$

$$H_c(\psi) = \frac{B_c(\psi) - B_{i-1}}{\mu_c} + H_{i-1}$$

$$H'_C(\psi) = \frac{1}{\mu_c}$$

Next, $H_c(\psi)$ and $H'_c(\psi)$ were plugged into $f(\psi)$ and $f'(\psi)$. Then, for Newton Raphson, ψ was updated at each iteration using

$$\psi_{i+1} = \psi_i - \frac{f_i}{f'_i}$$

The whole implementation can be seen in `non_linear.py`. Program took 3 iterations to reach $|f(\psi) / f(0)| < 10^{-6}$ and ψ was found as 0.00016126936944407854 Wb as displayed in Figure 1.5.1.

```
bash-3.2$ cd /Users/dafneculha/Desktop/A3-260785524 ; /usr/bin/env /Library/Frameworks/Python.framework/Versions/3.9/bin/python3 /Users/dafneculha/.vscode/extensions/m
s-python.python-2020.11.371526539/pythonFiles/Lib/python/debugpy/launcher 51279 -- /Users/dafneculha/Desktop/A3-260785524/non_linear.py

Solving with Newton-Raphson
Initial flux = 0

Iteration # 1
Flux = 0.00019995383179510732

Iteration # 2
Flux = 0.00016892691694374336

Iteration # 3
Flux = 0.00016126936944407854
```

Figure 1.5.1: Solution of non-linear equation using Newton-Raphson

(f) Try solving the same problem with successive substitution. If the method does not converge, suggest and test a modification of the method that does converge.

For successive substitution, ψ was updated at each iteration using

$$\psi_{i+1} = \psi_i - f_i$$

Using this, with the method didn't converge, and the program kept running forever. Thus, program was modified using the steps explained below.

$$\mathcal{F} = N \cdot I = R_{Total} \times \psi$$

$$8000 = \psi \cdot \left(\frac{H_C(\psi)}{\psi} \cdot L_C + R_A \right)$$

From this, $H_C(\psi)$ can be derived as

$$H_C(\psi) = \frac{8000 - R_A \cdot \psi}{L_C} = \frac{8000 - 39\,788\,735.77 \cdot \psi}{0.3}$$

Using Table 1 values, $B_C(\psi)$ was obtained as

$$\mu_c = \frac{\Delta B_c}{\Delta H_c} = \frac{B_{i+1} - B_i}{H_{i+1} - H_i}$$

$$B_c(\psi) = \mu_c(H_i(\psi) - H_{i-1}) + B_{i-1}$$

And using $B_C(\psi)$ values, ψ was obtained as

$$\psi = Area \cdot B_c(\psi)$$

These values were plugged into $f(\psi)$

$$f(\psi) = 0 = H_C(\psi) \times 0.3 + 39\,788\,735.77 \cdot \psi - 8000$$

and successive substitution was run again. The whole implementation can be seen in `non_linear.py`

New program took 15 iterations to reach $|f(\psi) / f(0)| < 10^{-6}$ and ψ was found as 0.00016126939623324393 Wb as displayed in Figure 1.6.1.


```

Solving with Successive Substitution
Initial flux = 0

Iteration # 1
Flux = 0.00019460292539069513

Iteration # 2
Flux = 0.00013605231845920413

Iteration # 3
Flux = 0.00016983297272672076

Iteration # 4
Flux = 0.00015740236647371832

Iteration # 5
Flux = 0.00016258258720894833

Iteration # 6
Flux = 0.00016082340627539523

Iteration # 7
Flux = 0.00016142081663977025

Iteration # 8
Flux = 0.00016121793862042868

Iteration # 9
Flux = 0.00016128683513308797

Iteration # 10
Flux = 0.00016126343817036122

Iteration # 11
Flux = 0.00016127138367941107

Iteration # 12
Flux = 0.00016126868541832485

Iteration # 13
Flux = 0.00016126960173631577

Iteration # 14
Flux = 0.00016126929055862256

Iteration # 15
Flux = 0.00016126939623324393

```

Figure 1.6.1: Solution of non-linear equation using Successive Substitution

2. For the circuit shown in Figure 2.1.1 below, the DC voltage E is 200 mV, the resistance R is $512\ \Omega$, the reverse saturation current for diode A is $I_{SA} = 0.8\ \mu\text{A}$, the reverse saturation current for diode B is $I_{SB} = 1.1\ \mu\text{A}$ and assume $kT/q = 25\ \text{mV}$.

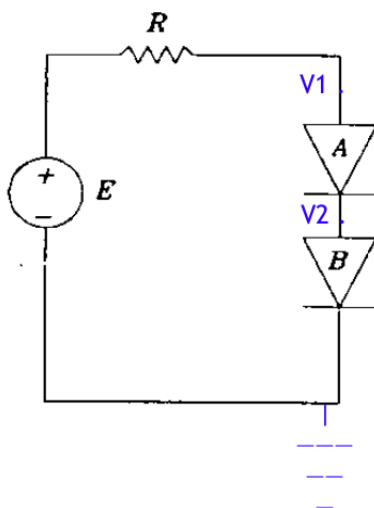


Figure 2.1.1: Circuit with Non-Linear Elements (Diodes)

(a) Derive nonlinear equations for a vector of nodal voltages, v_n , in the form $f(v_n) = 0$. Give f explicitly in terms of the variables I_{sA} , I_{sB} , E , R and v_n .

By Ohm's Law,

$$I = \frac{E - V1}{R}$$

By diode equations learnt in Microelectronics,

$$Vt = \frac{k \cdot T}{q} = 0.025V$$

KCL in Diode A,

$$I = I_{sA} \times \left[\exp\left(\frac{V1 - V2}{Vt}\right) - 1 \right] = I_{sA} \times \left[\exp\left(\frac{V1 - V2}{0.025}\right) - 1 \right]$$

$$\frac{E - V1}{R} = I_{sA} \times \left[\exp\left(\frac{V1 - V2}{0.025}\right) - 1 \right]$$

$$E - V1 = R \times I_{sA} \times \left[\exp\left(\frac{V1 - V2}{0.025}\right) - 1 \right]$$

$$f(V1) = 0 = V1 - E + R \times I_{sA} \times \left[\exp\left(\frac{V1 - V2}{0.025}\right) - 1 \right]$$

KCL in Diode B,

$$I = I_{sB} \times \left[\exp\left(\frac{V2}{Vt}\right) - 1 \right] = I_{sB} \times \left[\exp\left(\frac{V2}{0.025}\right) - 1 \right]$$

$$\frac{E - V1}{R} = I_{sB} \times \left[\exp\left(\frac{V2}{0.025}\right) - 1 \right] = I_{sA} \times \left[\exp\left(\frac{V1 - V2}{Vt}\right) - 1 \right]$$

$$f(V2) = I_{sA} \times \left[\exp\left(\frac{V1 - V2}{Vt}\right) - 1 \right] - I_{sB} \times \left[\exp\left(\frac{V2}{0.025}\right) - 1 \right]$$

$$f(Vn) = 0 = \begin{bmatrix} f(V1) \\ f(V2) \end{bmatrix} = \begin{bmatrix} V1 - E + R \times I_{sA} \times \left[\exp\left(\frac{V1 - V2}{0.025}\right) - 1 \right] \\ I_{sA} \times \left[\exp\left(\frac{V1 - V2}{0.025}\right) - 1 \right] - I_{sB} \times \left[\exp\left(\frac{V2}{0.025}\right) - 1 \right] \end{bmatrix}$$

(b) Solve the equation $f = 0$ by the Newton-Raphson method. At each step, record f and the voltage across each diode. Is the convergence quadratic? [Hint: define a suitable error measure ϵ_k].

To design this program, $f(V_n)$ found in previous part was used. Partial derivatives of f_1 and f_2 were taken with respect to V_1 and V_2 were taken and they were used to define the Jacobi matrix as presented below.

$$J = \begin{bmatrix} 1 + R \times I_{sA} \times \frac{1}{0.025} \times \left[\exp\left(\frac{V_1 - V_2}{0.025}\right) \right] & -R \times I_{sA} \times \frac{1}{0.025} \times \left[\exp\left(\frac{V_1 - V_2}{0.025}\right) \right] \\ I_{sA} \times \frac{1}{0.025} \times \left[\exp\left(\frac{V_1 - V_2}{0.025}\right) \right] & -I_{sA} \times \frac{1}{0.025} \times \left[\exp\left(\frac{V_1 - V_2}{0.025}\right) \right] - I_{sB} \times \frac{1}{0.025} \times \left[\exp\left(\frac{V_2}{0.025}\right) \right] \end{bmatrix}$$

And for each step $V_n = \begin{bmatrix} V_1 \\ V_2 \end{bmatrix}$ was updated as $V_{n+1} = \frac{-f + J \cdot V}{J} = -f * J^{-1} + V_n$

And $f(V_n)$ was updated with new V_n entries.

Initial guess for V_1 and V_2 were set as 0 volts. Since f_1 and f_2 are supposed to be 0, absolute value of them give the error. Error measure was set as 10^{-15} so that the program would keep iterating until f_1 and f_2 both become as small as 10^{-15} . Using these, f_1 , f_2 , V_1 and V_2 was computed after each iteration. The full implementation can be seen in Appendix at diodes.py. Results obtained are presented in Figure 2.2.1. It takes 6 iterations for the program to reach a tolerable result. V_1 was found as 0.18213962349633997 volts and V_2 was found as 0.08719379125672429 volts.

To see if the convergence is quadratic, f_1 and f_2 are observed. Looking at Figure 2.2.1, it can be observed that the number of zeros after decimal point in f_1 is 1, 2, 3, 7, 12, 'inf' as iterations increment from 1 to 6. f_2 operates in a similar fashion where the number of zeros after decimal point in f_2 is 0, 5, 7, 9, 14, 21 as iterations increment from 0 to 6. This confirms the expectation that the convergence is quadratic.

```
bash-3.2$ cd "/Users/dafneculha/Google Drive/School/Current Courses/ECSE 543/A3-26
0785524" ; /usr/bin/env /Library/Frameworks/Python.framework/Versions/3.9/bin/pytho
n3 /Users/dafneculha/.vscode/extensions/ms-python.python-2020.11.371526539/pythonFi
les/lib/python/debugpy/launcher 53649 -- "/Users/dafneculha/Google Drive/School/Cur
rent Courses/ECSE 543/A3-260785524/diodes.py"
```

```
Initial guesses
```

```
V1 = 0
```

```
V2 = 0
```

```
Initial residual
```

```
f1 = 0.0
```

```
f2 = -0.2
```

```
Iteration # 1
```

```
V1 = 0.19812073101961689
```

```
V2 = 0.08341925516615446
```

```
f1 = 0.03797622055434545
```

```
f2 = 4.8001808035091884e-05
```

```
Iteration # 2
```

```
V1 = 0.18413995441486622
```

```
V2 = 0.08433689566842323
```

```
f1 = 0.005918277697945149
```

```
f2 = 1.1538035340957757e-05
```

```
Iteration # 3
```

```
V1 = 0.18230748657784016
```

```
V2 = 0.08710806942596115
```

```
f1 = 0.0003541257681694866
```

```
f2 = 4.869660430844707e-07
```

```
Iteration # 4
```

```
V1 = 0.18214000800517066
```

```
V2 = 0.08719341483543555
```

```
f1 = 9.406043097598404e-07
```

```
f2 = 1.627918842005767e-09
```

```
Iteration # 5
```

```
V1 = 0.1821396235002092
```

```
V2 = 0.08719379125430769
```

```
f1 = 8.462917866491892e-12
```

```
f2 = 1.2450340780047198e-14
```

```
Iteration # 6
```

```
V1 = 0.18213962349633997
```

```
V2 = 0.08719379125672429
```

```
f1 = 0.0
```

```
f2 = -6.776263578034403e-21
```

```
bash-3.2$ █
```

Figure 2.2.1: V1, V2, f1 and f2 in each iteration

Appendix

1 - interpolation.py

```
from sympy import *
```

```
def lagrange_interpolation (points_x, points_y):
```

```
    param = symbols('B')
```

```
    n = len(points_x)
```

```
    yx = 0
```

```

Ljx = [0.0 for j in range (n)]
# For j = 0,1,2,...,n
for j in range (n):
    Fjx = 1
    Fjxj = 1
    for r in range (n):
        if (r != j):
            # Fjx = PRODUCT (x-xr)
            Fjx = Fjx * (param - points_x[r])
            Fjxj = Fjxj * (points_x[j] - points_x[r])
    #Ljx = Fjx / Fjxj
    Ljx[j] = Fjx.expand()/Fjxj
# yx = SUM (aj*Ljx)
for j in range (n):
    yx = yx + points_y[j] * Ljx[j]
return yx

```

```

def cubic_hermite(points_x,points_y):
    n = len(points_x)
    param = symbols('B')
    Ujx = []
    Vjx = []
    yx = 0
    # Lj(x)
    Ljx = [None for j in range (n)]
    #L'(x)
    Ljx_prime = [None for j in range (n)]
    b = [None for j in range (n)]
    a = [None for j in range (n)]

```

```

for j in range(n):
    Fjx = 1
    Fjxj = 1
    for r in range (n):
        if (r != j):
            # Fjx = PRODUCT (x-xr)
            Fjx = Fjx * (param - points_x[r])

```

```

    Fjxj = Fjxj * (points_x[j] - points_x[r])
# To obtain Lj(x)
# Lj(x) = Fjx / Fjxj
Ljx = Fjx.expand() / Fjxj

#To obtain L'j(x), diff Lj(x) / dx
Ljx_prime = lambdify(param, diff(Ljx))

U = (1 - 2 * Ljx_prime(points_x[j]) * (param - points_x[j]))*(Ljx**2)
Ujx.append(U)
V = (param - points_x[j])*(Ljx * Ljx)
Vjx.append(V)

# a[j] = y(j)
for j in range(n):
    a[j] = points_y[j]

# b[j] = y'(j)
for j in range(n):
    if (j < (n-1)):
        b[j] = (points_y[j + 1] - points_y[j]) / (points_x[j + 1] - points_x[j])
    elif (j == (n-1)):
        b[j] = points_y[j]/points_x[j]
    #a[j] = points_y[j]

# y(x) = SUM (a*Ujx + b*Vjx)
for j in range(n):
    yx = yx + a[j]*Ujx[j] + b[j]*Vjx[j]
yx = expand(yx)
return yx

if __name__ == "__main__":

    B = [0.0, 0.2, 0.4, 0.6, 0.8, 1.0, 1.1, 1.2, 1.3, 1.4, 1.5, 1.6, 1.7, 1.8, 1.9]
    H = [0.0, 14.7, 36.5, 71.7, 121.4, 197.4, 256.2, 348.7, 540.6, 1062.8, 2318.0, 4781.9, 8687.4, 13924.3, 22650.2]

```

```

print ("\nPart A\n")
points_x = [0.0, 0.2, 0.4, 0.6, 0.8, 1.0]
print ('B = ', points_x)
points_y = [0.0, 14.7, 36.5, 71.7, 121.4, 197.4]
print ('H = ', points_y)
a = lagrange_interpolation(points_x, points_y)
print ('H = ',a)

print ("\n")

print ('Part B\n')
points_x = [0.0, 1.3, 1.4, 1.7, 1.8, 1.9]
print ('B = ', points_x)
points_y = [0.0, 540.6, 1062.8, 8687.4, 13924.3, 22650.2]
print ('H = ', points_y)
b = lagrange_interpolation(points_x, points_y)
print ('H = ',b)

print ("\n")

print ('Part C\n')
points_x = [0.0, 1.3, 1.4, 1.7, 1.8, 1.9]
print ('B = ', points_x)
points_y = [0.0, 540.6, 1062.8, 8687.4, 13924.3, 22650.2]
print ('H = ', points_y)
c = cubic_hermite(points_x, points_y)
print ('H = ',c)

print ("\n")

```

2 - matrices.py

```

import math
import numpy as np

```

```
import time
import random
```

```
# returns True if matrix A is symmetric
```

```
def is_symmetric(A):
    n = len(A)
    if (n==1):
        return True
    for i in range(n):
        for j in range(i + 1, n):
            return A[i][j] == A[j][i]
#A = [[0.2]]
#print(is_symmetric(A))
```

```
# transforms col vectors to right format.
```

```
# [x,y,z] into [[x],[y],[z]]
```

```
def transform(A):
    try:
        cols = len(A[0])
        return A
    except TypeError:
        result = [[0] for a in range(len(A))]
        for i in range(0, len(A)):
            result[i][0] = A[i]
        return result
#A = [2,3,4]
#print(A)
#print(transform(A))
```

```
#returns the transpose of matrix A
```

```
def transpose(A):
    A = transform (A)
    A_transpose = [[0 for i in A] for j in A[0]]
    for i in range(len(A)):
        for j in range(len(A[0])):
            A_transpose[j][i] = A[i][j]
```



```

    return A_transpose
#A = [[2,3,0]]
#print(transpose(A))
#print(np.transpose(A))

#returns A+B
def add(A, B):
    A = transform(A)
    B = transform(B)
    if len(A)!= len(B) or len(A[0]) != len(B[0]):
        exit('addition not successful because of matrix size error')
    result = []
    for i in range(len(A)):
        result.append([A[i][k]+B[i][k] for k in range(len(A[0]))])
    return result
#A = [2,3,0]
#B = [1,3,0]
#print(add(A,B))

#returns A-B
def subtract(A, B):
    A = transform(A)
    B = transform(B)
    if len(A)!= len(B) or len(A[0]) != len(B[0]):
        exit('subtraction not successful because of matrix size error')
    result = []
    for i in range(len(A)):
        result.append([A[i][k]-B[i][k] for k in range(len(A[0]))])
    return result
#A = [[2,3,0], [2,2,1]]
#B = [[1,3,0], [5,1,1]]
#print(subtract(A,B))

#returns A*B
def dot_product(A, B):
    A = transform(A)
    B = transform(B)

```

```

if len(A[0])!= len(B):
    exit('dot product not successful because of matrix size error')
result = []
result = [[0 for i in range(len(B[0]))]for k in range(len(A))]
for i in range(len(A)):
    for j in range(len(B[0])):
        for k in range(len(A[0])):
            result[i][j] += A[i][k]*B[k][j]
    return result
#A = [[1,3], [0,1]]
#B = [3,1]
#print(dot_product(A,B))

```

#returns scalar*A

```

def scalar_product(scalar, A):
    A = transform(A)
    result = [[0 for i in range(len(A[0]))]for k in range(len(A))]
    for i in range(len(A)):
        for j in range(len(A[0])):
            result[i][j] = scalar*A[i][j]
    return result
#A = [2,6]
#print(scalar_product(5,A))

```

#returns determinant of a 2x2 matrix

```

def determinant(A):
    det = A[0][0] * A[1][1] - A[0][1] * A[1][0]
    return det

```

#returns product inverse of A

```

def inverse (A):
    det_A = determinant(A)
    inv_A = [[None for x in range (len(A))] for y in range(len(A))]
    inv_A[0][0] = A[1][1] / det_A
    inv_A[0][1] = -1 * A[0][1] / det_A
    inv_A[1][0] = -1 * A[1][0] / det_A
    inv_A[1][1] = A[0][0] / det_A

```

```

    return inv_A
#A = [[2,6],[3,1]]
#print(determinant(A))
#print (inverse(A))

#returns a random symmetric positive definite square matrix of size n
def random_symmetric_positive_definite_matrix(n):
    A = np.random.randint(-10,10, size=(n,n))
    return dot_product(A,transpose(A))

#returns a random symmetric positive definite vector of size n
def random_vector(n):
    return np.random.randint(-10,10, size=(n))

```

3 - diodes.py

```

from matrices import *
import numpy as np
import math

# constants in SI units
E = 0.2
R = 512
IsA = 0.8 * 10**(-6)
IsB = 1.1 * 10**(-6)
Vt = 0.025
tolerance = 10**(-15)

def newton_raphson (V1, V2, k):

    #initial residuals
    f1 = V1 - E + R * IsA * (math.exp((V1 - V2) / Vt) - 1)
    f2 = IsA * ((math.exp((V1 - V2) / Vt) - 1)) - IsB * (math.exp(V2 / Vt) - 1)

    print ("\nIteration #", k)

```

#Partial Derivatives

```
d_f1_V1 = 1.0 + (R * IsA / Vt) * (math.exp((V1 - V2) / Vt))
d_f1_V2 = -1.0 * (R * IsA / Vt) * (math.exp((V1 - V2) / Vt))
d_f2_V1 = (IsA / Vt) * (math.exp((V1 - V2) / Vt))
d_f2_V2 = -1.0 * (IsA / Vt) * (math.exp((V1 - V2) / Vt)) - (IsB / Vt) * (math.exp(V2 / Vt))
```

Jacobian Matrix

```
J = [[d_f1_V1, d_f1_V2], [d_f2_V1, d_f2_V2]]
```

```
V = [V1, V2]
```

```
f = [f1, f2]
```

#Update voltages

```
V = add(scalar_product(-1, dot_product(inverse(J), f)), V)
V1 = V[0][0]
V2 = V[1][0]
print ('V1 = ', V1)
print ('V2 = ', V2)
```

#Update residuals

```
f1 = V1 - E + R * IsA * (math.exp((V1 - V2) / Vt) - 1)
f2 = IsA * ((math.exp((V1 - V2) / Vt) - 1)) - IsB * (math.exp(V2 / Vt) - 1)

print ('f1 = ', f1)
print ('f2 = ', f2)
```

```
return V1, V2, f1, f2
```

```
if __name__ == "__main__":
```

initial guesses

```
V1 = 0
```

```
V2 = 0
```

#initial residuals

```
f1 = IsA * (math.exp((V1 - V2) / Vt) - 1) - IsB * (math.exp(V2 / Vt) - 1)
f2 = V1 - E + R * IsB * (math.exp(V2 / Vt) - 1)
```

```

print ("\nInitial guesses")
print("V1 = ", V1)
print("V2 = ", V2)
print ("\nInitial residual")
print('f1 = ', f1)
print('f2 = ', f2)

#counter
k = 1
while (abs(f1) >=tolerance or abs(f2)>=tolerance):
    V1, V2, f1, f2 = newton_raphson(V1, V2,k)
    k = k + 1
print ("\n")

```

4 – non_linear.py

```

import math
import numpy as np

B = [0.0, 0.2, 0.4, 0.6, 0.8, 1.0, 1.1, 1.2, 1.3, 1.4, 1.5, 1.6, 1.7, 1.8, 1.9]
H = [0.0, 14.7, 36.5, 71.7, 121.4, 197.4, 256.2, 348.7, 540.6, 1062.8, 2318.0, 4781.9, 8687.4, 13924.3, 22650.2]

Area = 10**(-4)
Lc = 0.3
La = 0.005
N = 800
I = 10
perm0 = 4*math.pi*10**(-7)

Ra = La / (perm0 * Area)
#print (Ra)
MMF = N * I

tolerance = 10**(-6)

# B = flux / area

```

```

def get_B (flux):
    B = flux / Area
    return B
#print(get_B(0.3))

# B = mu * H
# mu = B / H
# H = B / mu
# H_prime = 1/mu
def get_H_and_H_prime(flux):
    Bc = flux/Area
    if Bc > B[-1]:
        mu = (B[-1] - B[-2]) / (H[-1] - H[-2])
        Hc = (Bc - B[-1]) / mu + H[-1]
        H_prime = 1/mu
        return Hc, H_prime

    for i in range(len(B)):
        if B[i] > Bc:
            mu = (B[i] - B[i-1]) / (H[i] - H[i-1])
            Hc = (Bc - B[i-1]) / mu + H[i-1]
            H_prime = 1/mu
            return Hc, H_prime

def get_f(flux):
    Hc, H_prime = get_H_and_H_prime(flux)
    f = Hc * Lc + Ra * flux - MMF
    return f

def get_f_prime(flux):
    Hc, H_prime = get_H_and_H_prime(flux)
    f_prime = H_prime * Lc / Area + Ra
    return f_prime

def newton_raphson(flux):
    print ('Initial flux = ', flux)
    i = 1

```

```

while abs(get_f(flux)/get_f(0)) > tolerance:
    print ("\nIteration #", i)
    flux = flux - get_f(flux)/get_f_prime(flux)
    print ('Flux = ', flux)
    i += 1
return flux

```

```

def successive_sub_xxx(flux):
    i = 1
    while abs(get_f(flux)/get_f(0)) >= tolerance:
        print ("\nIteration #", i)
        flux = get_f(flux)
        print ('Flux = ', flux)
        i += 1
    return flux

```

```

def get_Bc(flux):
    Hc = (MMF - Ra * flux) / Lc
    for i in range(len(H) - 1):
        if H[i] <= Hc < H[i + 1]:
            mu = ((B[i + 1] - B[i])/(H[i + 1] - H[i]))
            Bc = B[i] + ((Hc) - H[i]) * mu
            return Bc
        elif Hc > H[- 1]:
            mu = ((B[- 1] - B[- 2])/(H[- 1] - H[- 2]))
            Bc = B[- 1] + (Hc - H[- 1]) * mu
            return Bc

```

```

def successive_sub(flux):
    i = 1
    print ('Initial flux = ', flux)
    while abs(get_f(flux)/get_f(0)) >= tolerance:
        print ("\nIteration #", i)
        flux = Area * get_Bc(flux)
        print ('Flux = ', flux)
        i += 1
    return flux

```

```
if __name__ == "__main__":  
    print ("\nSolving with Newton-Raphson")  
    newton_raphson(0)  
    print ("\n")  
  
    print ("\nSolving with Successive Substitution")  
    successive_sub (0)  
    print("\n")
```