HW3_1

November 21, 2018

```
In [311]: import numpy as np
          import pandas as pd
          f=open('Soft_Magnetic_Dataset.csv','r')
          data=f.read()
          f.close
Out[311]: <function TextIOWrapper.close()>
In [312]: def list_concat(word):
              out=''
              for i in word:
                   out+=i
              return(out)
In [313]: temp=[]
          data1=[]
          for t in data:
              if (t!='\n' \text{ and } t!=','):
                   temp.append(t)
              else:
                   data1.append(temp)
                   temp=[]
In [314]: data2=[]
          for i in data1:
              data2.append(list_concat(i))
In [315]: cols=data2[0:42]
          data3=data2[42:]
          print(len(data3)/42)
1294.0
In []:
In [316]: data4=np.array(data3)
          data5=np.reshape(data4,(1294,42))
          m,n=data5.shape
          import pdb
          #pdb.set_trace()
```

```
In [317]: data_df=pd.DataFrame(data=data5,columns=cols)
          data_df.head()
          data_df.iloc[0,-5] == ''
          data_df.head()
          print(data_df.columns)
Index(['Fe', 'Si', 'C', 'Al', 'B', 'P', 'Ga', 'Ge', 'Cu', 'Ag', 'Au', 'Zn',
       'Ti', 'V', 'Cr', 'Zr', 'Nb', 'Mo', 'Hf', 'Ta', 'W', 'Ce', 'Pr', 'Gd',
       'U', 'Annealing temperature (K)', 'Annealing Time (s)',
       'Primary Crystallization Onset (K)',
       'Primary Crystallization Peak (K) ',
       'Secondary Crystallization Peak (K)', ' Longitudinal Annealing field',
       'Transverse Annealing field', 'Ribbon Thickness (um) ',
       'Coercivity (A/m)', 'Curie Temperature (K)',
       'Core Loss (kW/m<sup>3</sup> @ 0.2T 100kHz)',
       'Electrical Resistivity (10^-6*ohm*m)', 'Permeability @ 1kHz',
       'Magnetostriction (10^-6)', 'Magnetic Saturation (T)',
       'Grain Diameter (nm)', 'Reference DOI'],
      dtype='object')
In [318]: #preprocessing
          #remove columns with zero elements
          nonzero_ratio=np.zeros(len(cols))
          counter=0
          for col_i in cols:
              n_zeros=0
              #print('column : ',col_i)
              #import pdb; pdb.set_trace()
              for j in range(0,len(data_df[col_i])):
                  if(data_df[col_i][j]=='0' or data_df[col_i][j]==''):
                      n_zeros+=1
              nonzero_ratio[counter]=1-n_zeros/len(data_df[col_i])
              counter+=1
          print('non zero ratio : ',nonzero_ratio)
          drop_cols=[]
          for i in range(0,len(data_df.columns)-1):
              if (nonzero_ratio[i]<.05):</pre>
                  #print('non zero ratio', nonzero_ratio[i])
                  drop_cols.append(cols[i])
          #add the last column
          drop_cols.append(cols[-1])
          data_df_cleancols=data_df.drop(drop_cols,axis=1)
          new_cols=data_df_cleancols.columns
```

```
print(len(new_cols),len(cols))
         #import pdb; pdb.set_trace()
         data_df_cleancols.head()
         print(data_df_cleancols.columns)
                 [1.00000000e+00 7.34157651e-01 3.94126739e-02 9.65996909e-02
non zero ratio :
 9.76043277e-01 1.69242658e-01 2.08655332e-02 7.34157651e-02
 7.18701700e-01 1.23647604e-02 2.55023184e-02 7.72797527e-04
 6.95517774e-03 1.93199382e-02 2.39567233e-02 1.01236476e-01
 5.83462133e-01 8.50077280e-02 7.72797527e-03 3.55486862e-02
 5.79598145e-02 2.31839258e-03 7.72797527e-04 1.54559505e-03
 6.18238022e-03 1.00000000e+00 1.00000000e+00 8.88717156e-02
 4.08809892e-01 3.50850077e-01 5.40958269e-03 7.72797527e-03
 7.94435858e-01 5.72642968e-01 7.26429675e-02 2.93663060e-02
 3.70942813e-02 2.86707883e-01 1.56105100e-01 2.27202473e-01
 1.57650696e-01 1.00000000e+00]
23 42
Index(['Fe', 'Si', 'Al', 'B', 'P', 'Ge', 'Cu', 'Zr', 'Nb', 'Mo', 'W',
       'Annealing temperature (K)', 'Annealing Time (s)',
       'Primary Crystallization Onset (K)',
       'Primary Crystallization Peak (K) ',
       'Secondary Crystallization Peak (K)', 'Ribbon Thickness (um) ',
       'Coercivity (A/m)', 'Curie Temperature (K)', 'Permeability @ 1kHz',
       'Magnetostriction (10^-6)', 'Magnetic Saturation (T)',
       'Grain Diameter (nm)'],
      dtype='object')
In [319]: drop_rows=[]
         coer=data_df_cleancols['Coercivity (A/m)']
         for i in range(0, len(coer)):
             if coer[i]=='':
                 drop_rows.append(i)
         print('Rows to be dropped',len(drop_rows))
         data_df_cleanrows=data_df_cleancols.drop(drop_rows)
         print('Remaining Rows',len(data_df_cleanrows))
         data_df_cleanrows=data_df_cleanrows.reset_index(drop=True)
         data_df_cleanrows.head()
Rows to be dropped 553
Remaining Rows 741
Out[319]:
              Fe
                    Si Al B P Ge Cu Zr Nb Mo
         0 72.5 13.5 0 9 0 1 1 0 3 0
         1 71.5
                  9.5 1 9 0 5 1 0 3 0
         2 73.5 13.5 0 9 0 0 1 0 3 0
         3 73.5 13.5 0 9 0 0 1 0 3 0
         4 73.5 13.5 0 7 0 2 1 0 3 0
```

```
0
          1
          2
          3
          4
            Secondary Crystallization Peak (K) Ribbon Thickness (um) Coercivity (A/m)
                                                                              4644.936707
                                                                     25
                                                                              4106.876281
          1
          2
                                                                     25
                                                                              3922.927079
          3
                                                                     25
                                                                              2428.806915
          4
                                                                     25
                                                                              996.4925244
            Curie Temperature (K) Permeability @ 1kHz Magnetostriction (10^-6) \
          0
          1
          2
          3
          4
            Magnetic Saturation (T) Grain Diameter (nm)
          1
          2
          3
          4
          [5 rows x 23 columns]
In [320]: np.random.seed(0)
          for col in new_cols:
              if col != 'Coercivity (A/m)':
                  for i in range(0,len(data_df_cleanrows)):
                      try:
                           data_df_cleanrows[col][i]=float(data_df_cleanrows[col][i])+np.random.n
                      except ValueError:
                               pass
          data_df_cleanrows.head()
          print(data_df_cleanrows.columns)
Index(['Fe', 'Si', 'Al', 'B', 'P', 'Ge', 'Cu', 'Zr', 'Nb', 'Mo', 'W',
       'Annealing temperature (K)', 'Annealing Time (s)',
       \hbox{'Primary Crystallization Onset (K)',}\\
       'Primary Crystallization Peak (K) ',
       'Secondary Crystallization Peak (K)', 'Ribbon Thickness (um)',
       'Coercivity (A/m)', 'Curie Temperature (K)', 'Permeability @ 1kHz',
```

Primary Crystallization Onset (K) Primary Crystallization Peak (K)

```
'Magnetostriction (10^-6)', 'Magnetic Saturation (T)',
       'Grain Diameter (nm)'],
      dtype='object')
In [321]: m,n=data_df_cleanrows.shape
          for col in new_cols:
              if col!='Coercivity (A/m)':
                  for j in range(0,len(data_df_cleanrows)):
                      try:
                          data_df_cleanrows[col][j]=data_df_cleanrows[col][j][0]
                      except ValueError:
                          pass
                      except IndexError:
                          pass
                      except TypeError:
                          pass
In [322]: m,n=data_df_cleanrows.shape
          for i in range(m):
              for j in range(n):
                  if (data_df_cleanrows.iloc[i][j]==''):
                      data_df_cleanrows.iloc[i][j]=np.nan
In [323]: data_df_cleanrows.head()
Out[323]:
                  Fe
                           Si
                                      Al
                                                В
                                                          Ρ
                                                                  Ge
                                                                             Cu \
                                -2.40747 8.15242 0.625268
            76.0281
                      11.2196
                                                            1.37394
                                                                      -1.11413
          1 72.3003
                        13.01
                                 1.83409 7.80666 -0.386928
                                                             2.88894
                                                                      -2.64939
                                          10.1341
          2 75.4575
                       13.234
                                 1.37576
                                                     2.4826
                                                             1.19122
                                                                       1.48348
          3 77.9818 11.9686 0.0997145
                                         10.9765 -0.311796
                                                            -2.7506 0.709696
          4 77.2351 14.6116
                                 2.69607 5.97219 -1.47834 3.24602
                                                                       2.00399
                   Zr
                             Nb
                                      Mo
            -2.08559
                        5.99051 1.37661
          1 -1.31121
                        2.4048 1.32479
          2 0.514308
                        1.50777
                                 1.77779
          3 -0.713619
                        3.64489 1.08875
            -4.50449 0.183463 -1.45304
            Primary Crystallization Onset (K) Primary Crystallization Peak (K)
          0
                                          NaN
                                                                             NaN
          1
                                          NaN
                                                                             NaN
          2
                                          NaN
                                                                             NaN
          3
                                          NaN
                                                                             NaN
          4
                                          NaN
                                                                             NaN
            Secondary Crystallization Peak (K) Ribbon Thickness (um) Coercivity (A/m)
          0
                                           NaN
                                                              22.8576
                                                                            4644.936707
```

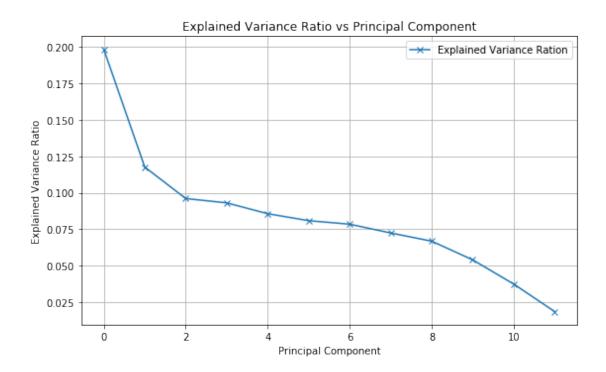
```
NaN
                                                                 26.1108
                                                                               4106.876281
          1
          2
                                                                 21.5819
                                             NaN
                                                                               3922.927079
          3
                                             NaN
                                                                 23.6256
                                                                               2428.806915
          4
                                             NaN
                                                                 25.1247
                                                                               996.4925244
            Curie Temperature (K) Permeability @ 1kHz Magnetostriction (10^-6)
          0
                               NaN
                                                    NaN
          1
                               NaN
                                                    NaN
                                                                               NaN
          2
                               NaN
                                                    NaN
                                                                               NaN
          3
                               NaN
                                                    NaN
                                                                               NaN
          4
                               NaN
                                                    NaN
                                                                               NaN
            Magnetic Saturation (T) Grain Diameter (nm)
          0
                                 NaN
                                                      NaN
          1
                                 NaN
                                                      NaN
          2
                                 NaN
                                                      NaN
          3
                                 NaN
                                                      NaN
                                 NaN
                                                      NaN
          [5 rows x 23 columns]
In [324]: from sklearn.preprocessing import StandardScaler
          scaler=StandardScaler()
          for col in new_cols:
              if col != 'Coercivity (A/m)':
                  null_values=data_df_cleanrows[col].isnull()
                  data_df_cleanrows.loc[~null_values,[col]]=scaler.fit_transform(data_df_cleanro
In [325]: data_df_cleanrows.head()
Out [325]:
                     Fe
                               Si
                                           Al
                                                       В
                                                                  Ρ
                                                                           Ge
                                                                                       Cu
             -0.215755 0.624717
                                    -1.17265 -0.0661397 -0.272111
                                                                    0.222183
                                                                               -0.913517
          1
            -0.818957
                        0.909259
                                    0.680074
                                               -0.173938 -0.550679
                                                                      0.72874
                                                                                -1.65492
              -0.30809
                         0.944867
                                    0.479876
                                                0.551689
                                                            0.23905
                                                                     0.161089
                                                                                 0.340914
              0.100374
                        0.743754 -0.0775066
                                                0.814312 -0.530002
                                                                     -1.15691 -0.0327589
                                               -0.745866 -0.851049
          4 -0.0204459
                           1.1638
                                      1.05659
                                                                     0.848136
                                                                                 0.592277
                    Zr
                              Nb
                                        Мо
                                                                  \
          0 -0.984052
                         1.66608 0.534391
          1 -0.677158  0.320822  0.509423
          2 0.046319 -0.015721
                                  0.727698
          3 -0.440323 0.786069
                                  0.395685
                                                     . . .
          4 -1.94269 -0.512567
                                  -0.82906
                                                    . . .
            Primary Crystallization Onset (K) Primary Crystallization Peak (K)
          0
                                            NaN
                                                                                NaN
          1
                                            NaN
                                                                                NaN
          2
                                            NaN
                                                                                NaN
```

```
3
                                            NaN
                                                                                NaN
          4
                                                                                NaN
                                            NaN
            Secondary Crystallization Peak (K) Ribbon Thickness (um) Coercivity (A/m)
          0
                                                               0.0261461
                                                                               4644.936707
                                             NaN
          1
                                             NaN
                                                                0.576264
                                                                               4106.876281
          2
                                             NaN
                                                                -0.189586
                                                                               3922.927079
          3
                                             NaN
                                                                 0.156016
                                                                               2428.806915
          4
                                             NaN
                                                                0.409522
                                                                               996.4925244
            Curie Temperature (K) Permeability @ 1kHz Magnetostriction (10^-6) \
          0
                               NaN
                                                     NaN
                                                                               NaN
          1
                                                     NaN
                               NaN
                                                                               {\tt NaN}
          2
                                NaN
                                                     NaN
                                                                               NaN
          3
                                NaN
                                                     NaN
                                                                               {\tt NaN}
          4
                               NaN
                                                     NaN
                                                                               NaN
            Magnetic Saturation (T) Grain Diameter (nm)
          0
                                  NaN
                                                       NaN
          1
                                  NaN
                                                       NaN
          2
                                  NaN
                                                       NaN
          3
                                  NaN
                                                       NaN
                                  NaN
                                                       NaN
          [5 rows x 23 columns]
In [334]: """Obtain the features to apply PCA on"""
          new_cols1=[]
          cols=[]
          for col in new_cols:
               cols.append(col)
          #print(type(cols),cols)
          new_cols1=cols[0:12]
          print(new_cols1,len(new_cols1))
['Fe', 'Si', 'Al', 'B', 'P', 'Ge', 'Cu', 'Zr', 'Nb', 'Mo', 'W', 'Annealing temperature (K)'] 12
In [336]: from sklearn.decomposition import PCA
          pca = PCA(n_components=12)
          print(data_df_cleanrows[new_cols1].isnull().sum().sum())
          data=data_df_cleanrows[new_cols1]
          X=data[data.notna()]
          pca.fit(X)
          print(X.isnull().sum().sum())
0
```

0

```
In [333]: #find number of NAN values
          for i in range(0,12):
              null_values=data_df_cleanrows[new_cols[i]].isnull()
              type(null_values)
              print(len(data_df_cleanrows.loc[~null_values,[new_cols[i]]]))
          print(new_cols[0:12])
741
741
741
741
741
741
741
741
741
741
741
741
Index(['Fe', 'Si', 'Al', 'B', 'P', 'Ge', 'Cu', 'Zr', 'Nb', 'Mo', 'W',
       'Annealing temperature (K)'],
      dtype='object')
In [367]:
Out [367]: 0.07339656151675539
In [343]: #plot error vs n
          import matplotlib.pyplot as plt
          fig, ax = plt.subplots(figsize=[8,5])
          plt.plot(pca.explained_variance_ratio_,marker='x',label='Explained Variance Ration')
          plt.hold(True)
          plt.title('Explained Variance Ratio vs Principal Component')
          plt.ylabel('Explained Variance Ratio')
          plt.xlabel('Principal Component')
          fig.tight_layout()
          ax.legend()
          plt.grid(True)
          plt.show
          fig.savefig('hw3_3a.png')
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:5: MatplotlibDeprecationWarning
    Future behavior will be consistent with the long-time default:
    plot commands add elements without first clearing the
    Axes and/or Figure.
C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\__init__.py:911: MatplotlibDeprecationWarn
  mplDeprecation)
```

C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\rcsetup.py:156: MatplotlibDeprecationWarni
mplDeprecation)

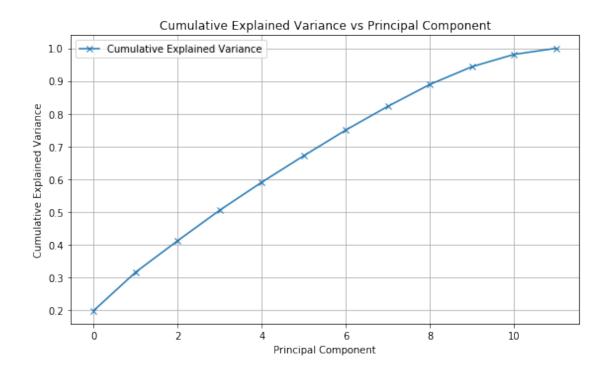


```
In [345]: cum_ratio=[]
          cum=0
          for ratio in pca.explained_variance_ratio_:
              cum+=ratio
              cum_ratio.append(cum)
          #plot error vs n
          fig, ax = plt.subplots(figsize=[8,5])
          plt.plot(cum_ratio,marker='x',label='Cumulative Explained Variance')
         plt.hold(True)
          plt.title('Cumulative Explained Variance vs Principal Component')
          plt.ylabel('Cumulative Explained Variance')
          plt.xlabel('Principal Component')
          fig.tight_layout()
          ax.legend()
          plt.grid(True)
         plt.show
          fig.savefig('hw3_3b.png')
```

C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:10: MatplotlibDeprecationWarnin Future behavior will be consistent with the long-time default: plot commands add elements without first clearing the

Axes and/or Figure.

- # Remove the CWD from sys.path while we load stuff.
- C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib__init__.py:911: MatplotlibDeprecationWarn
 mplDeprecation)
- C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\rcsetup.py:156: MatplotlibDeprecationWarni
 mplDeprecation)



PC1=np.array(pca.components_[0])
PC2=np.array(pca.components_[1])
PC3=np.array(pca.components_[2])

b=np.array(X.iloc[740])

np.dot(a,b)

```
if data_df_cleanrows[new_cols[-6]][i]==0:
                  plt.plot(x,y,'r',marker='x')
                  plt.hold(True)
              elif data_df_cleanrows[new_cols[-6]][i]==1:
                  plt.plot(x,y,'g',marker='x')
                  plt.hold(True)
              else:
                  plt.plot(x,y,'b',marker='x')
                  plt.hold(True)
          plt.title('PC1 vs PC2')
          plt.ylabel('PC2')
          plt.xlabel('PC1')
          fig.tight_layout()
          ax.legend()
          plt.grid(True)
          plt.show
          fig.savefig('hw3_4a.png')
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:23: MatplotlibDeprecationWarnin
    Future behavior will be consistent with the long-time default:
    plot commands add elements without first clearing the
    Axes and/or Figure.
C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\__init__.py:911: MatplotlibDeprecationWarn
  mplDeprecation)
C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\rcsetup.py:156: MatplotlibDeprecationWarni
  mplDeprecation)
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:20: MatplotlibDeprecationWarnin
    Future behavior will be consistent with the long-time default:
    plot commands add elements without first clearing the
    Axes and/or Figure.
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:17: MatplotlibDeprecationWarnin
    Future behavior will be consistent with the long-time default:
    plot commands add elements without first clearing the
    Axes and/or Figure.
No handles with labels found to put in legend.
```

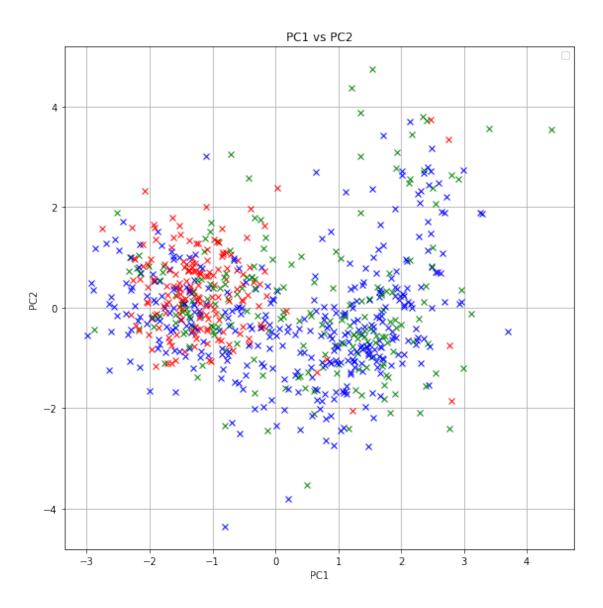
fig, ax = plt.subplots(figsize=[8,8])

pt=np.array(X.iloc[i])

x=np.dot(pt,PC1)
y=np.dot(pt,PC2)

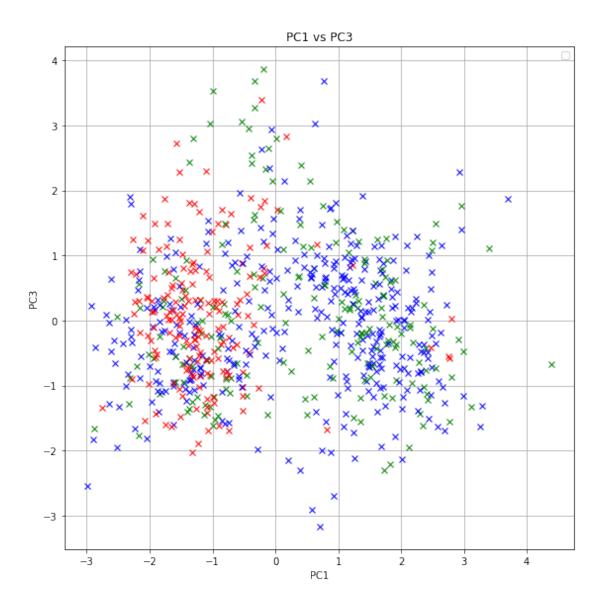
for i in range(0,len(data_df_cleanrows[new_cols[-6]])):

#first PC1 and PC2



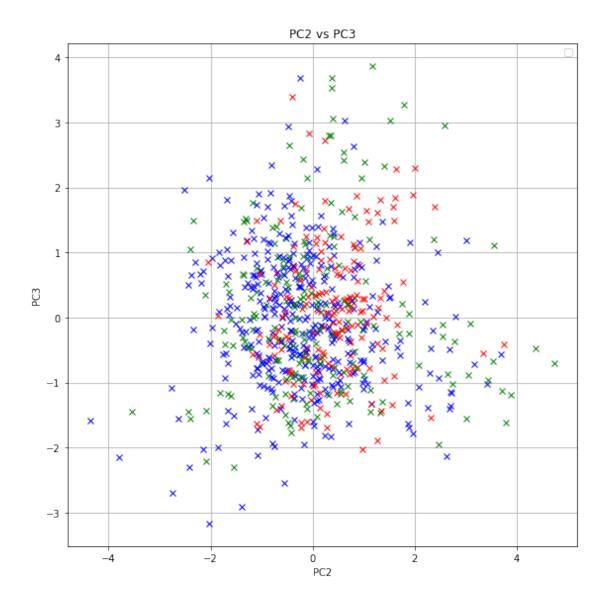
```
In [375]: fig, ax = plt.subplots(figsize=[8,8])
    #next PC1 and PC3
    for i in range(0,len(data_df_cleanrows[new_cols[-6]])):
        pt=np.array(X.iloc[i])
        x=np.dot(pt,PC1)
        y=np.dot(pt,PC3)
        if data_df_cleanrows[new_cols[-6]][i]==0:
            plt.plot(x,y,'r',marker='x')
            plt.hold(True)
        elif data_df_cleanrows[new_cols[-6]][i]==1:
            plt.plot(x,y,'g',marker='x')
            plt.hold(True)
        else:
```

```
plt.plot(x,y,'b',marker='x')
                plt.hold(True)
         plt.title('PC1 vs PC3')
         plt.ylabel('PC3')
         plt.xlabel('PC1')
         fig.tight_layout()
         ax.legend()
         plt.grid(True)
         plt.show
         fig.savefig('hw3_4b.png')
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:15: MatplotlibDeprecationWarnin
   Future behavior will be consistent with the long-time default:
   plot commands add elements without first clearing the
   Axes and/or Figure.
  from ipykernel import kernelapp as app
mplDeprecation)
C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\rcsetup.py:156: MatplotlibDeprecationWarni
 mplDeprecation)
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:12: MatplotlibDeprecationWarnin
   Future behavior will be consistent with the long-time default:
   plot commands add elements without first clearing the
   Axes and/or Figure.
  if sys.path[0] == '':
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:9: MatplotlibDeprecationWarning
   Future behavior will be consistent with the long-time default:
   plot commands add elements without first clearing the
   Axes and/or Figure.
  if __name__ == '__main__':
No handles with labels found to put in legend.
```



```
In [381]: fig, ax = plt.subplots(figsize=[8,8])
    #finally PC2 and PC23
    for i in range(0,len(data_df_cleanrows[new_cols[-6]])):
        pt=np.array(X.iloc[i])
        x=np.dot(pt,PC2)
        y=np.dot(pt,PC3)
        if data_df_cleanrows[new_cols[-6]][i]==0:
            plt.plot(x,y,'r',marker='x')
            plt.hold(True)
        elif data_df_cleanrows[new_cols[-6]][i]==1:
            plt.plot(x,y,'g',marker='x')
            plt.hold(True)
        else:
```

```
plt.plot(x,y,'b',marker='x')
                plt.hold(True)
         plt.title('PC2 vs PC3')
         plt.ylabel('PC3')
         plt.xlabel('PC2')
         fig.tight_layout()
         ax.legend()
         plt.grid(True)
         plt.show
         fig.savefig('hw3_4c.png')
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:15: MatplotlibDeprecationWarnin
   Future behavior will be consistent with the long-time default:
   plot commands add elements without first clearing the
   Axes and/or Figure.
  from ipykernel import kernelapp as app
mplDeprecation)
C:\Users\aksha\Anaconda3\lib\site-packages\matplotlib\rcsetup.py:156: MatplotlibDeprecationWarni
 mplDeprecation)
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:12: MatplotlibDeprecationWarnin
   Future behavior will be consistent with the long-time default:
   plot commands add elements without first clearing the
   Axes and/or Figure.
  if sys.path[0] == '':
C:\Users\aksha\Anaconda3\lib\site-packages\ipykernel_launcher.py:9: MatplotlibDeprecationWarning
   Future behavior will be consistent with the long-time default:
   plot commands add elements without first clearing the
   Axes and/or Figure.
  if __name__ == '__main__':
No handles with labels found to put in legend.
```



In [380]: print(pca.get_covariance())

```
[-0.03768452  0.06163139  -0.00307291  0.04944866  -0.03769337
                                                                                                           0.03965982
     1.00135135 0.00407828 0.00415234 -0.00591753 -0.04126098
                                                                                                         0.01776972]
  \hbox{ [ 0.18188581 -0.16066039 -0.02590294 -0.10655042 -0.09254025 -0.01132244 ] } 
     0.00407828 1.00135135 -0.06498361 -0.07867549 0.00425655 0.01936066]
 [-0.31582802  0.18883969  0.01138607  0.06350481  -0.1546759
                                                                                                           0.11318762
     0.00415234 -0.06498361 1.00135135 0.02141887 -0.05351544 0.1946975 ]
   \begin{bmatrix} -0.09259027 & 0.12074562 & -0.05249769 & 0.0483009 & -0.05103195 & -0.00814852 \end{bmatrix} 
   -0.00591753 -0.07867549 0.02141887 1.00135135 0.02497942 0.05115617
   \begin{bmatrix} 0.06143245 & -0.11500165 & -0.06988509 & 0.01772845 & -0.06564444 & -0.07006339 \end{bmatrix} 
   -0.04126098 0.00425655 -0.05351544 0.02497942 1.00135135 0.09330368]
 [-0.28264495 0.36640789 0.0164819
                                                                  0.01906001 -0.32328742 -0.05778079
                                                                  0.05115617 0.09330368 1.00135135]]
     0.01776972 0.01936066 0.1946975
In [387]: """The loading matrix W, we have the PCs in the columns and the original variables in
                 comp_mat=pca.components_
                 df_components=pd.DataFrame(np.transpose(comp_mat),columns=['PC1','PC2','PC3','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','PC4','
                 df_components
Out[387]:
                                                                  PC3
                                                                                   PC4
                                                                                                    PC5
                                                                                                                      PC6
                                PC1
                                                 PC2
                                                                                                                                       PC7 \
                       0.495333 - 0.332875 \quad 0.010266 \quad 0.009189 - 0.055634 - 0.016890 - 0.052244
                      -0.533270 0.111963 -0.156031 -0.050673 -0.174002 0.076273 -0.155193
                     -0.006857 0.461403 -0.177270 0.096178 -0.248446 -0.319845 0.234003
                     -0.185512 -0.468177 0.333770 -0.275425 -0.210867 -0.407294 0.046601
                       0.396329 \quad 0.477389 \ -0.109116 \ -0.157813 \quad 0.045252 \quad 0.099786 \quad 0.081968
                    -0.043539 0.158239 0.627001 0.211257 0.471186 0.127733
                                                                                                                            0.159545
                     -0.061899 -0.012510 0.302562 0.089112 -0.575218 0.532712 0.498176
                 7
                       0.131639 -0.231033 -0.099144  0.696356 -0.035059  0.220873 -0.294988
                     -0.306006 0.141902 0.197357 0.171336 0.325309 -0.148213 0.065714
                     -0.125895 -0.048304 -0.077698 -0.492704 0.220865
                                                                                                            0.580332 -0.248261
                 10 0.033875 -0.326175 -0.386266 -0.053671
                                                                                           0.383246
                                                                                                            0.031405 0.688125
                 11 -0.380837 -0.096149 -0.363489 0.273602 0.052764
                                                                                                            0.084433 0.097049
                                PC8
                                                 PC9
                                                                PC10
                                                                                                   PC12
                                                                                  PC11
                       0.020629 0.191008 -0.364859 -0.379433
                                                                                            0.568805
                      -0.162216 -0.237466 0.243980 -0.097995
                                                                                            0.681293
                       2
                                                                                            0.069030
                 3
                       -0.135396  0.165318  0.003406  0.636431  0.329803
                 5
                       0.177553 -0.402384 -0.179897 0.027607 0.212028
                 6
                     7
                       0.315756  0.062744  0.345906  0.260491  0.079861
                      -0.093367  0.796974  0.120510  -0.129115  0.103576
                       0.058694 -0.083969 0.309676 -0.002294 0.112255
```

0.03965982 -0.01132244 0.11318762 -0.00814852 -0.07006339 -0.05778079]

```
In [392]: """Most important feature"""
          #if we look at PC1 then we can decide that feature 1 is the most important feature
          new_cols1[0:2]
Out[392]: ['Fe', 'Si']
In [405]: """Hierarchical Clustering"""
          from scipy.cluster.hierarchy import dendrogram, linkage
          X1=data_df_cleanrows[new_cols1[0:2]]
          m, n=X1.shape
         X = \Gamma I
          labels1=[]
          for i in range (0,132,12):
              X.append(X1.iloc[i])
              if data_df_cleanrows['Coercivity (A/m)'][i]==0:
                  labels1.append('Low')
              elif data_df_cleanrows['Coercivity (A/m)'][i]==1:
                  labels1.append('Medium')
              else:
                  labels1.append('High')
          #single linkage
          #import pdb;pdb.set_trace()
          Z=linkage(X,'single')
          fig, ax = plt.subplots(figsize=[8,8])
          dendrogram(Z,labels=labels1)
          plt.title('Dendrogram of samples with Fe and Si as Features- Single Linkage')
          plt.show()
          fig.savefig('hw3_6a.png')
          #complete
          Z=linkage(X,'complete')
          fig, ax = plt.subplots(figsize=[8,8])
          dendrogram(Z,labels=labels1)
          plt.title('Dendrogram of samples with Fe and Si as Features- Complete Linkage')
          plt.show()
          fig.savefig('hw3_6b.png')
          #average linkage
          Z=linkage(X,'average')
          fig, ax = plt.subplots(figsize=[8,8])
          dendrogram(Z,labels=labels1)
          plt.title('Dendrogram of samples with Fe and Si as Features- Average Linkage')
         plt.show()
          fig.savefig('hw3_6c.png')
```

