

ML0101EN-Clus-Hierarchical-Cars-py-v1

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Hierarchical Clustering

Welcome to Lab of Hierarchical Clustering with Python using Scipy and Scikit-learn package.

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Hierarchical Clustering - Agglomerative

We will be looking at a clustering technique, which is Agglomerative Hierarchical Clustering. Remember that agglomerative is the bottom up approach. In this lab, we will be looking at Agglomerative clustering, which is more popular than Divisive clustering. We will also be using Complete Linkage as the Linkage Criteria. NOTE: You can also try using Average Linkage wherever Complete Linkage would be used to see the difference!

```
[1]: import numpy as np
import pandas as pd
from scipy import ndimage
from scipy.cluster import hierarchy
from scipy.spatial import distance_matrix
from matplotlib import pyplot as plt
from sklearn import manifold, datasets
from sklearn.cluster import AgglomerativeClustering
from sklearn.datasets.samples_generator import make_blobs
%matplotlib inline
```

Generating Random Data

We will be generating a set of data using the `make_blobs` class. Input these parameters into `make_blobs`:

`n_samples`: The total number of points equally divided among clusters.

Choose a number from 10-1500

`centers`: The number of centers to generate, or the fixed center locations.

Choose arrays of x,y coordinates for generating the centers. Have 1-10 centers (ex. `centers=[[1,1], [2,5]]`)

`cluster_std`: The standard deviation of the clusters. The larger the number, the further apart the clusters

Choose a number between 0.5-1.5

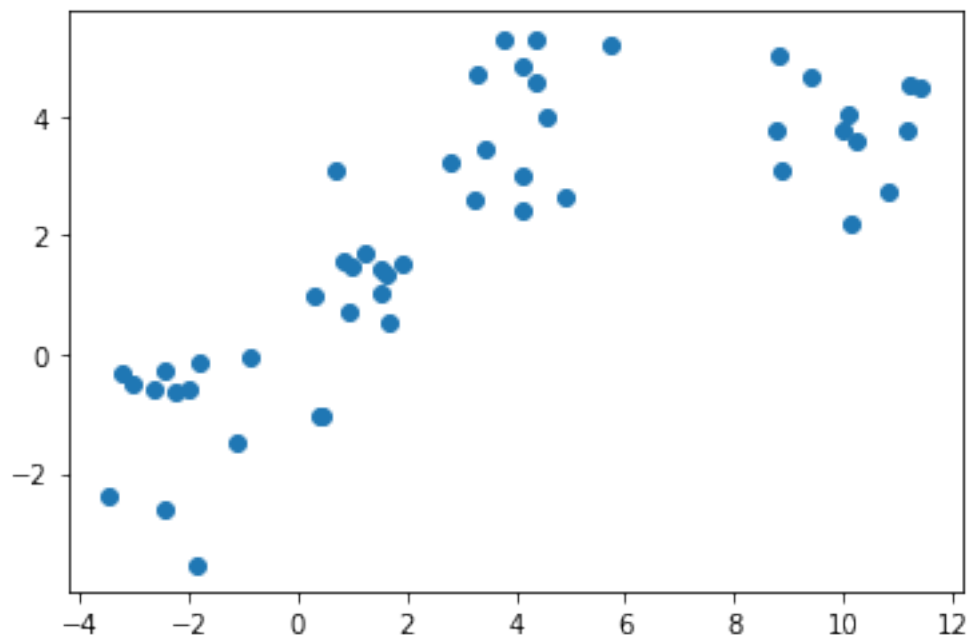
Save the result to `X1` and `y1`.

```
[2]: X1, y1 = make_blobs(n_samples=50, centers=[[4,4], [-2, -1], [1, 1], [10,4]], cluster_std=0.9)
```

Plot the scatter plot of the randomly generated data

```
[3]: plt.scatter(X1[:, 0], X1[:, 1], marker='o')
```

```
[3]: <matplotlib.collections.PathCollection at 0x7f6fb4bff908>
```



Agglomerative Clustering

We will start by clustering the random data points we just created.

The Agglomerative Clustering class will require two inputs:

`n_clusters`: The number of clusters to form as well as the number of centroids to generate.

Value will be: 4

`linkage`: Which linkage criterion to use. The linkage criterion determines which distance to use between sets of observation. The algorithm will merge the pairs of cluster that minimize this criterion.

Value will be: 'complete'

Note: It is recommended you try everything with 'average' as well
Save the result to a variable called agglom

```
[4]: agglom = AgglomerativeClustering(n_clusters = 4, linkage = 'average')
```

Fit the model with X2 and y2 from the generated data above.

```
[5]: agglom.fit(X1,y1)
```

```
[5]: AgglomerativeClustering(affinity='euclidean', compute_full_tree='auto',  
    connectivity=None, linkage='average', memory=None,  
    n_clusters=4, pooling_func='deprecated')
```

Run the following code to show the clustering! Remember to read the code and comments to gain more understanding on how the plotting works.

```
[6]: # Create a figure of size 6 inches by 4 inches.  
plt.figure(figsize=(6,4))  
  
# These two lines of code are used to scale the data points down,  
# Or else the data points will be scattered very far apart.  
  
# Create a minimum and maximum range of X1.  
x_min, x_max = np.min(X1, axis=0), np.max(X1, axis=0)  
  
# Get the average distance for X1.  
X1 = (X1 - x_min) / (x_max - x_min)  
  
# This loop displays all of the datapoints.  
for i in range(X1.shape[0]):  
    # Replace the data points with their respective cluster value  
    # (ex. 0) and is color coded with a colormap (plt.cm.spectral)  
    plt.text(X1[i, 0], X1[i, 1], str(y1[i]),  
            color=plt.cm.nipy_spectral(agglom.labels_[i] / 10.),  
            fontdict={'weight': 'bold', 'size': 9})  
  
# Remove the x ticks, y ticks, x and y axis  
plt.xticks([])  
plt.yticks([])  
plt.axis('off')  
  
# Display the plot of the original data before clustering  
plt.scatter(X1[:, 0], X1[:, 1], marker='.')  
# Display the plot  
plt.show()
```



Dendrogram Associated for the Agglomerative Hierarchical Clustering

Remember that a distance matrix contains the distance from each point to every other point of a dataset. Use the function `distance_matrix`, which requires two inputs. Use the Feature Matrix, `X2` as both inputs and save the distance matrix to a variable called `dist_matrix`. Remember that the distance values are symmetric, with a diagonal of 0's. This is one way of making sure your matrix is correct. (print out `dist_matrix` to make sure it's correct)

```
[7]: dist_matrix = distance_matrix(X1,X1)
      print(dist_matrix)
```

```
[[0.         0.41529098 0.41121742 ... 0.49194307 0.47495725 0.06664725]
 [0.41529098 0.         0.05817215 ... 0.07705311 0.78406558 0.37248808]
 [0.41121742 0.05817215 0.         ... 0.09788586 0.74928598 0.36133842]
 ...
 [0.49194307 0.07705311 0.09788586 ... 0.         0.8470407 0.4476907 ]
 [0.47495725 0.78406558 0.74928598 ... 0.8470407 0.         0.46441301]
 [0.06664725 0.37248808 0.36133842 ... 0.4476907 0.46441301 0.         ]]
```

Using the linkage class from `hierarchy`, pass in the parameters:

The distance matrix

'complete' for complete linkage

Save the result to a variable called `Z`

```
[8]: Z = hierarchy.linkage(dist_matrix, 'complete')
```

```
/home/jupyterlab/conda/envs/python/lib/python3.6/site-
packages/ipynb_launcher.py:1: ClusterWarning: scipy.cluster: The symmetric
non-negative hollow observation matrix looks suspiciously like an uncondensed
distance matrix
```

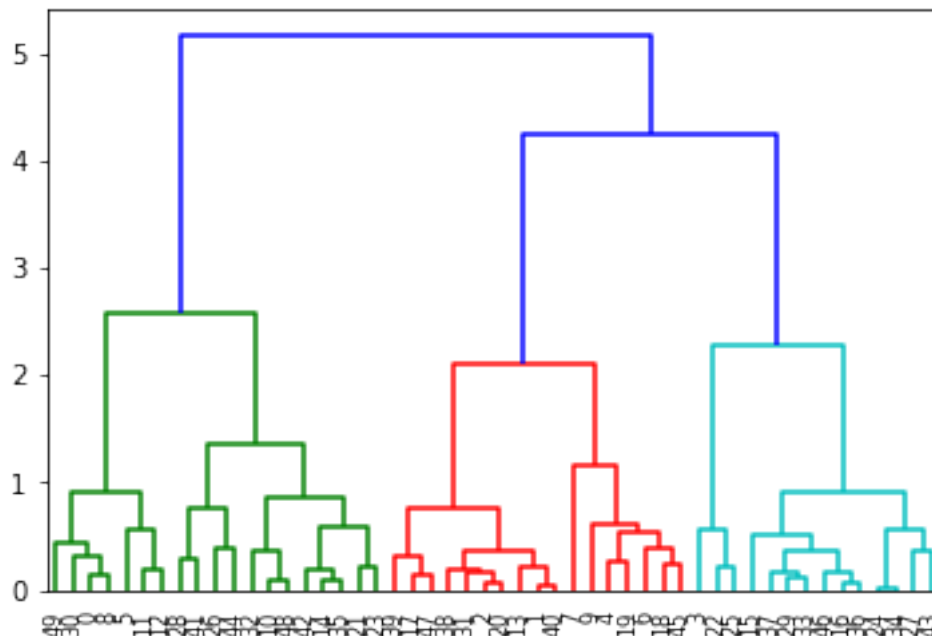
```
"""Entry point for launching an IPython kernel.
```

A Hierarchical clustering is typically visualized as a dendrogram as shown in the following cell. Each merge is represented by a horizontal line. The y-coordinate of the horizontal line is the similarity of the two clusters that were merged, where cities are viewed as singleton clusters. By moving up from the bottom layer to the top node, a dendrogram allows us to reconstruct the history of merges that resulted in the depicted clustering.

Next, we will save the dendrogram to a variable called `dendro`. In doing this, the dendrogram will also be displayed. Using the `dendrogram` class from `hierarchy`, pass in the parameter:

`Z`

```
[9]: dendro = hierarchy.dendrogram(Z)
```

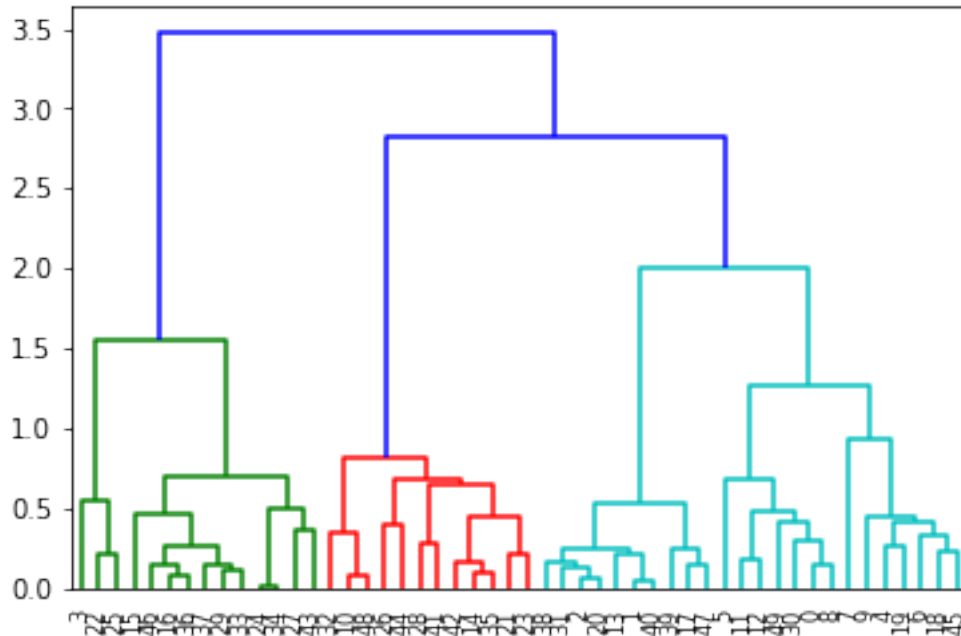


0.1 Practice

We used **complete** linkage for our case, change it to **average** linkage to see how the dendrogram changes.

```
[10]: # write your code here
Z = hierarchy.linkage(dist_matrix, 'average')
dendro = hierarchy.dendrogram(Z)
```

```
/home/jupyterlab/conda/envs/python/lib/python3.6/site-
packages/ipykernel_launcher.py:2: ClusterWarning: scipy.cluster: The symmetric
non-negative hollow observation matrix looks suspiciously like an uncondensed
distance matrix
```



Double-click **here** for the solution.

Clustering on Vehicle dataset

Imagine that an automobile manufacturer has developed prototypes for a new vehicle. Before introducing the new model into its range, the manufacturer wants to determine which existing vehicles on the market are most like the prototypes—that is, how vehicles can be grouped, which group is the most similar with the model, and therefore which models they will be competing against.

Our objective here, is to use clustering methods, to find the most distinctive clusters of vehicles. It will summarize the existing vehicles and help manufacturers to make decision about the supply of new models.

0.1.1 Download data

To download the data, we will use `!wget` to download it from IBM Object Storage.

Did you know? When it comes to Machine Learning, you will likely be working with large datasets. As a business, where can you host your data? IBM is offering a unique opportunity for businesses, with 10 Tb of IBM Cloud Object Storage: [Sign up now for free](#)

```
[11]: !wget -O cars_clus.csv https://s3-api.us-geo.objectstorage.softlayer.net/cf-courses-data/
      ↪ CognitiveClass/ML0101ENv3/labs/cars_clus.csv
```

```
--2019-08-15 01:18:19-- https://s3-api.us-geo.objectstorage.softlayer.net/cf-
courses-data/CognitiveClass/ML0101ENv3/labs/cars_clus.csv
Resolving s3-api.us-geo.objectstorage.softlayer.net (s3-api.us-
geo.objectstorage.softlayer.net)... 67.228.254.193
Connecting to s3-api.us-geo.objectstorage.softlayer.net (s3-api.us-
geo.objectstorage.softlayer.net)|67.228.254.193|:443... connected.
```

HTTP request sent, awaiting response... 200 OK
Length: 17774 (17K) [text/csv]
Saving to: cars_clus.csv

cars_clus.csv 100%[=====>] 17.36K --.-KB/s in 0.1s

2019-08-15 01:18:20 (168 KB/s) - cars_clus.csv saved [17774/17774]

0.2 Read data

lets read dataset to see what features the manufacturer has collected about the existing models.

```
[12]: filename = 'cars_clus.csv'

#Read csv
pdf = pd.read_csv(filename)
print ("Shape of dataset: ", pdf.shape)

pdf.head(5)
```

Shape of dataset: (159, 16)

```
[12]:  manufact  model  sales  resale  type  price  engine_s  horsepower  wheelbas  \
0   Acura  Integra  16.919  16.360  0.000  21.500    1.800   140.000   101.200
1   Acura    TL    39.384  19.875  0.000  28.400    3.200   225.000   108.100
2   Acura    CL   14.114  18.225  0.000   $null$    3.200   225.000   106.900
3   Acura    RL    8.588  29.725  0.000  42.000    3.500   210.000   114.600
4   Audi    A4   20.397  22.255  0.000  23.990    1.800   150.000   102.600

      width  length  curb_wgt  fuel_cap   mpg  lnsales  partition
0  67.300  172.400   2.639   13.200  28.000   2.828         0.0
1  70.300  192.900   3.517   17.200  25.000   3.673         0.0
2  70.600  192.000   3.470   17.200  26.000   2.647         0.0
3  71.400  196.600   3.850   18.000  22.000   2.150         0.0
4  68.200  178.000   2.998   16.400  27.000   3.015         0.0
```

The feature sets include price in thousands (price), engine size (engine_s), horsepower (horsepower), wheelbase (wheelbas), width (width), length (length), curb weight (curb_wgt), fuel capacity (fuel_cap) and fuel efficiency (mpg).

Data Cleaning

lets simply clear the dataset by dropping the rows that have null value:

```
[13]: print ("Shape of dataset before cleaning: ", pdf.size)
pdf[[ 'sales', 'resale', 'type', 'price', 'engine_s',
      'horsepow', 'wheelbas', 'width', 'length', 'curb_wgt', 'fuel_cap',
      'mpg', 'lnsales']] = pdf[['sales', 'resale', 'type', 'price', 'engine_s',
      'horsepow', 'wheelbas', 'width', 'length', 'curb_wgt', 'fuel_cap',
      'mpg', 'lnsales']].apply(pd.to_numeric, errors='coerce')
pdf = pdf.dropna()
```

```
pdf = pdf.reset_index(drop=True)
print("Shape of dataset after cleaning: ", pdf.size)
pdf.head(5)
```

Shape of dataset before cleaning: 2544

Shape of dataset after cleaning: 1872

```
[13]:  manufact  model  sales  resale  type  price  engine_s  horsepower \
0   Acura  Integra  16.919  16.360   0.0  21.50      1.8      140.0
1   Acura    TL    39.384  19.875   0.0  28.40      3.2      225.0
2   Acura    RL    8.588  29.725   0.0  42.00      3.5      210.0
3   Audi    A4    20.397  22.255   0.0  23.99      1.8      150.0
4   Audi    A6    18.780  23.555   0.0  33.95      2.8      200.0

      wheelbas  width  length  curb_wgt  fuel_cap  mpg  lnsales  partition
0      101.2    67.3   172.4    2.639    13.2  28.0    2.828        0.0
1      108.1    70.3   192.9    3.517    17.2  25.0    3.673        0.0
2      114.6    71.4   196.6    3.850    18.0  22.0    2.150        0.0
3      102.6    68.2   178.0    2.998    16.4  27.0    3.015        0.0
4      108.7    76.1   192.0    3.561    18.5  22.0    2.933        0.0
```

0.2.1 Feature selection

Lets select our feature set:

```
[14]: featureset = pdf[['engine_s', 'horsepow', 'wheelbas', 'width', 'length', 'curb_wgt', 'fuel_cap', 'mpg']]
```

0.2.2 Normalization

Now we can normalize the feature set. **MinMaxScaler** transforms features by scaling each feature to a given range. It is by default (0, 1). That is, this estimator scales and translates each feature individually such that it is between zero and one.

```
[15]: from sklearn.preprocessing import MinMaxScaler
x = featureset.values #returns a numpy array
min_max_scaler = MinMaxScaler()
feature_mtx = min_max_scaler.fit_transform(x)
feature_mtx [0:5]

array([[0.11428571, 0.21518987, 0.18655098, 0.28143713, 0.30625832,
        0.2310559 , 0.13364055, 0.43333333],
       [0.31428571, 0.43037975, 0.3362256 , 0.46107784, 0.5792277 ,
        0.50372671, 0.31797235, 0.33333333],
       [0.35714286, 0.39240506, 0.47722343, 0.52694611, 0.62849534,
        0.60714286, 0.35483871, 0.23333333],
       [0.11428571, 0.24050633, 0.21691974, 0.33532934, 0.38082557,
        0.34254658, 0.28110599, 0.4       ],
       [0.25714286, 0.36708861, 0.34924078, 0.80838323, 0.56724368,
```



```
0.5173913 , 0.37788018, 0.23333333]])
```

Clustering using Scipy

In this part we use Scipy package to cluster the dataset:

First, we calculate the distance matrix.

```
[16]: import scipy
      leng = feature_mtx.shape[0]
      D = scipy.zeros([leng,leng])
      for i in range(leng):
          for j in range(leng):
              D[i,j] = scipy.spatial.distance.euclidean(feature_mtx[i], feature_mtx[j])
```

In agglomerative clustering, at each iteration, the algorithm must update the distance matrix to reflect the distance of the newly formed cluster with the remaining clusters in the forest. The following methods are supported in Scipy for calculating the distance between the newly formed cluster and each: - single - complete - average - weighted - centroid

We use **complete** for our case, but feel free to change it to see how the results change.

```
[17]: import pylab
      import scipy.cluster.hierarchy
      Z = hierarchy.linkage(D, 'complete')
```

```
/home/jupyterlab/conda/envs/python/lib/python3.6/site-
packages/ipykernel_launcher.py:3: ClusterWarning: scipy.cluster: The symmetric
non-negative hollow observation matrix looks suspiciously like an uncondensed
distance matrix
```

This is separate from the ipykernel package so we can avoid doing imports until

Essentially, Hierarchical clustering does not require a pre-specified number of clusters. However, in some applications we want a partition of disjoint clusters just as in flat clustering. So you can use a cutting line:

```
[18]: from scipy.cluster.hierarchy import fcluster
      max_d = 3
      clusters = fcluster(Z, max_d, criterion='distance')
      clusters
```

```
[18]: array([ 1,  5,  5,  6,  5,  4,  6,  5,  5,  5,  5,  5,  4,  4,  5,  1,  6,
          5,  5,  5,  4,  2, 11,  6,  6,  5,  6,  5,  1,  6,  6, 10,  9,  8,
          9,  3,  5,  1,  7,  6,  5,  3,  5,  3,  8,  7,  9,  2,  6,  6,  5,
          4,  2,  1,  6,  5,  2,  7,  5,  5,  5,  4,  4,  3,  2,  6,  6,  5,
          7,  4,  7,  6,  6,  5,  3,  5,  5,  6,  5,  4,  4,  1,  6,  5,  5,
          5,  6,  4,  5,  4,  1,  6,  5,  6,  6,  5,  5,  5,  7,  7,  7,  2,
          2,  1,  2,  6,  5,  1,  1,  1,  7,  8,  1,  1,  6,  1,  1],
      dtype=int32)
```

Also, you can determine the number of clusters directly:

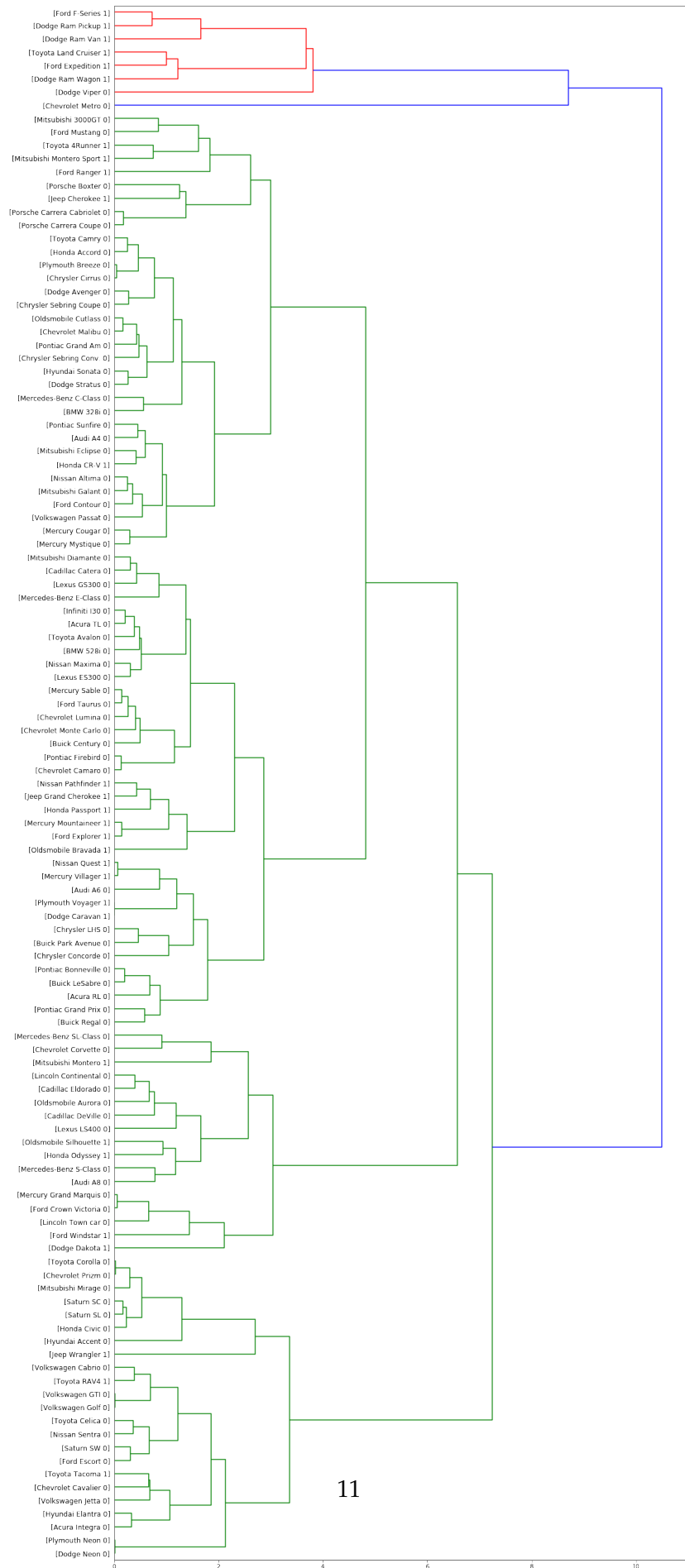
```
[19]: from scipy.cluster.hierarchy import fcluster
      k = 5
      clusters = fcluster(Z, k, criterion='maxclust')
```

```
clusters
```

```
[19]: array([1, 3, 3, 3, 3, 2, 3, 3, 3, 3, 3, 3, 2, 2, 3, 1, 3, 3, 3, 3, 2, 1,  
          5, 3, 3, 3, 3, 3, 1, 3, 3, 4, 4, 4, 4, 2, 3, 1, 3, 3, 3, 2, 3, 2,  
          4, 3, 4, 1, 3, 3, 3, 2, 1, 1, 3, 3, 1, 3, 3, 3, 3, 2, 2, 2, 1, 3,  
          3, 3, 3, 2, 3, 3, 3, 3, 2, 3, 3, 3, 3, 2, 2, 1, 3, 3, 3, 3, 3, 2,  
          3, 2, 1, 3, 3, 3, 3, 3, 3, 3, 3, 3, 3, 1, 1, 1, 1, 3, 3, 1, 1, 1,  
          3, 4, 1, 1, 3, 1, 1], dtype=int32)
```

Now, plot the dendrogram:

```
[20]: fig = pylab.figure(figsize=(18,50))  
def llf(id):  
    return '%s %s %s' % (pdf['manufact'][id], pdf['model'][id], int(float(pdf['type'][id])))  
  
dendro = hierarchy.dendrogram(Z, leaf_label_func=llf, leaf_rotation=0, leaf_font_size =12,  
    ↳orientation = 'right')
```



Clustering using scikit-learn

Lets redo it again, but this time using scikit-learn package:

```
[22]: dist_matrix = distance_matrix(feature_mtx, feature_mtx)
      print(dist_matrix)
```

```
[[0.          0.57777143 0.75455727 ... 0.28530295 0.24917241 0.18879995]
 [0.57777143 0.          0.22798938 ... 0.36087756 0.66346677 0.62201282]
 [0.75455727 0.22798938 0.          ... 0.51727787 0.81786095 0.77930119]
 ...
 [0.28530295 0.36087756 0.51727787 ... 0.          0.41797928 0.35720492]
 [0.24917241 0.66346677 0.81786095 ... 0.41797928 0.          0.15212198]
 [0.18879995 0.62201282 0.77930119 ... 0.35720492 0.15212198 0.          ]]
```

Now, we can use the 'AgglomerativeClustering' function from scikit-learn library to cluster the dataset. The AgglomerativeClustering performs a hierarchical clustering using a bottom up approach. The linkage criteria determines the metric used for the merge strategy:

- Ward minimizes the sum of squared differences within all clusters. It is a variance-minimizing approach and in this sense is similar to the k-means objective function but tackled with an agglomerative hierarchical approach.
- Maximum or complete linkage minimizes the maximum distance between observations of pairs of clusters.
- Average linkage minimizes the average of the distances between all observations of pairs of clusters.

```
[23]: agglom = AgglomerativeClustering(n_clusters = 6, linkage = 'complete')
      agglom.fit(feature_mtx)
      agglom.labels_
```

```
[23]: array([1, 2, 2, 1, 2, 3, 1, 2, 2, 2, 2, 2, 3, 3, 2, 1, 1, 2, 2, 2, 5, 1,
           4, 1, 1, 2, 1, 2, 1, 1, 1, 5, 0, 0, 0, 3, 2, 1, 2, 1, 2, 3, 2, 3,
           0, 3, 0, 1, 1, 1, 2, 3, 1, 1, 1, 2, 1, 1, 2, 2, 2, 3, 3, 3, 1, 1,
           1, 2, 1, 2, 2, 1, 1, 2, 3, 2, 3, 1, 2, 3, 5, 1, 1, 2, 3, 2, 1, 3,
           2, 3, 1, 1, 2, 1, 1, 2, 2, 2, 1, 1, 1, 1, 1, 1, 1, 1, 2, 1, 1, 1,
           2, 0, 1, 1, 1, 1, 1])
```

And, we can add a new field to our dataframe to show the cluster of each row:

```
[24]: pdf['cluster_'] = agglom.labels_
      pdf.head()
```

```
[24]:  manufact  model  sales  resale  type  price  engine_s  horsepower \
0   Acura  Integra  16.919  16.360   0.0  21.50         1.8       140.0
1   Acura    TL    39.384  19.875   0.0  28.40         3.2       225.0
2   Acura    RL    8.588  29.725   0.0  42.00         3.5       210.0
3   Audi    A4   20.397  22.255   0.0  23.99         1.8       150.0
4   Audi    A6   18.780  23.555   0.0  33.95         2.8       200.0
```

	wheelbas	width	length	curb_wgt	fuel_cap	mpg	lnsales	partition \
0	101.2	67.3	172.4	2.639	13.2 28.0	2.828	0.0	
1	108.1	70.3	192.9	3.517	17.2 25.0	3.673	0.0	
2	114.6	71.4	196.6	3.850	18.0 22.0	2.150	0.0	
3	102.6	68.2	178.0	2.998	16.4 27.0	3.015	0.0	
4	108.7	76.1	192.0	3.561	18.5 22.0	2.933	0.0	

	cluster_
0	1
1	2
2	2
3	1
4	2

```
[25]: import matplotlib.cm as cm
n_clusters = max(agglom.labels_)+1
colors = cm.rainbow(np.linspace(0, 1, n_clusters))
cluster_labels = list(range(0, n_clusters))

# Create a figure of size 6 inches by 4 inches.
plt.figure(figsize=(16,14))

for color, label in zip(colors, cluster_labels):
    subset = pdf[pdf.cluster_ == label]
    for i in subset.index:
        plt.text(subset.horsepow[i], subset.mpg[i],str(subset['model'][i]), rotation=25)
    plt.scatter(subset.horsepow, subset.mpg, s= subset.price*10, c=color,
        →label='cluster'+str(label),alpha=0.5)
# plt.scatter(subset.horsepow, subset.mpg)
plt.legend()
plt.title('Clusters')
plt.xlabel('horsepow')
plt.ylabel('mpg')
```

'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points.

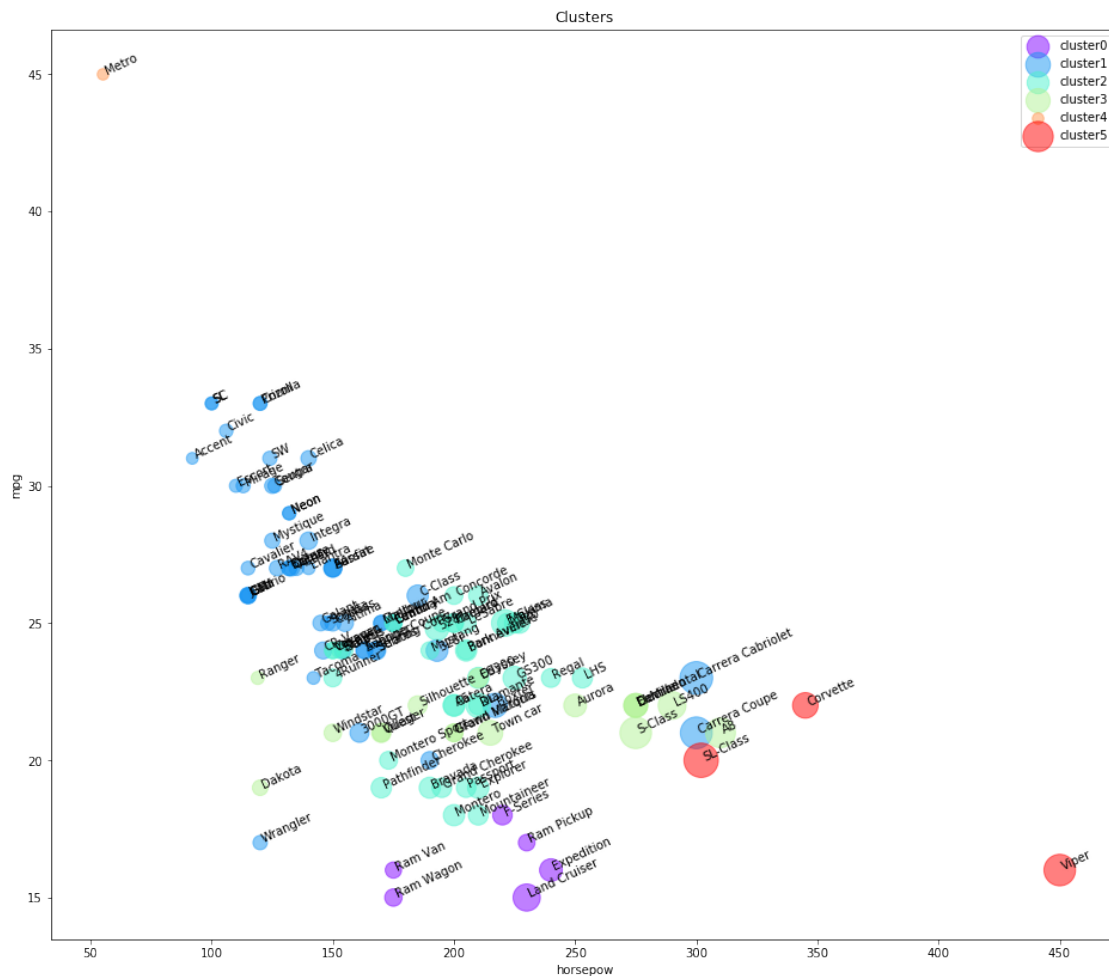
'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points.

'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points.

'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with

'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points.

[25]: Text(0, 0.5, 'mpg')



As you can see, we are seeing the distribution of each cluster using the scatter plot, but it is not very clear where is the centroid of each cluster. Moreover, there are 2 types of vehicles in our dataset, “truck” (value of 1 in the type column) and “car” (value of 1 in the type column). So, we use them to distinguish the classes, and summarize the cluster. First we count the number of cases in each group:

```
[26]: pdf.groupby(['cluster_', 'type'])['cluster_'].count()
```

```
[26]: cluster_  type
0         1.0    6
1         0.0   47
         1.0    5
2         0.0   27
         1.0   11
3         0.0   10
         1.0    7
4         0.0    1
5         0.0    3
Name: cluster_, dtype: int64
```

Now we can look at the characteristics of each cluster:

```
[27]: agg_cars = pdf.groupby(['cluster_', 'type'])['horsepow', 'engine_s', 'mpg', 'price'].mean()
agg_cars
```

```
[27]:      horsepow  engine_s      mpg      price
cluster_ type
0         1.0  211.666667  4.483333  16.166667  29.024667
1         0.0  146.531915  2.246809  27.021277  20.306128
         1.0  145.000000  2.580000  22.200000  17.009200
2         0.0  203.111111  3.303704  24.214815  27.750593
         1.0  182.090909  3.345455  20.181818  26.265364
3         0.0  256.500000  4.410000  21.500000  42.870400
         1.0  160.571429  3.071429  21.428571  21.527714
4         0.0   55.000000  1.000000  45.000000   9.235000
5         0.0  365.666667  6.233333  19.333333  66.010000
```

It is obvious that we have 3 main clusters with the majority of vehicles in those.

Cars: - Cluster 1: with almost high mpg, and low in horsepower. - Cluster 2: with good mpg and horsepower, but higher price than average. - Cluster 3: with low mpg, high horsepower, highest price.

Trucks: - Cluster 1: with almost highest mpg among trucks, and lowest in horsepower and price. - Cluster 2: with almost low mpg and medium horsepower, but higher price than average. - Cluster 3: with good mpg and horsepower, low price.

Please notice that we did not use **type**, and **price** of cars in the clustering process, but Hierarchical clustering could forge the clusters and discriminate them with quite high accuracy.

```
[28]: plt.figure(figsize=(16,10))
for color, label in zip(colors, cluster_labels):
    subset = agg_cars.loc[(label,)]
    for i in subset.index:
        plt.text(subset.loc[i][0]+5, subset.loc[i][2], 'type='+str(int(i)) + ', price='+str(int(subset.
        ↳loc[i][3]))+'k')
    plt.scatter(subset.horsepow, subset.mpg, s=subset.price*20, c=color,
    ↳label='cluster'+str(label))
plt.legend()
plt.title('Clusters')
```

```
plt.xlabel('horsepow')  
plt.ylabel('mpg')
```

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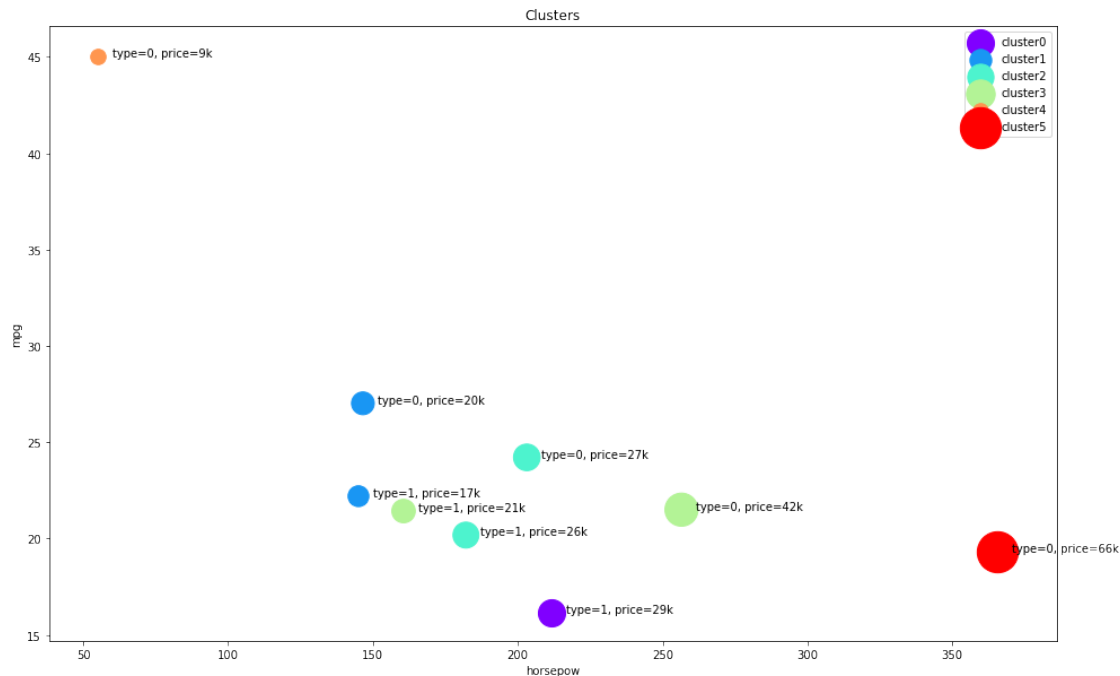
'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points.

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[28]: Text(0, 0.5, 'mpg')



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Thanks for completing this lesson!

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Saeed Aghabozorgi, PhD is a Data Scientist in IBM with a track record of developing enterprise level applications that substantially increases clients' ability to turn data into actionable knowledge. He is a researcher in data mining field and expert in developing advanced analytic methods like machine learning and statistical modelling on large datasets.

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