

# T81-558: Applications of Deep Neural Networks

#### Module 8: Kaggle Data Sets

- Instructor: Jeff Heaton, McKelvey School of Engineering, Washington University in St. Louis
- For more information visit the class website.

#### Module 8 Material

- Part 8.1: Introduction to Kaggle [Video] [Notebook]
- Part 8.2: Building Ensembles with Scikit-Learn and Keras [Video] [Notebook]
- Part 8.3: How Should you Architect Your Keras Neural Network: Hyperparameters [Video] [Notebook]
- Part 8.4: Bayesian Hyperparameter Optimization for Keras [Video] [Notebook]
- Part 8.5: Current Semester's Kaggle [Video] [Notebook]

## Part 8.1: Introduction to Kaggle

Kaggle runs competitions where data scientists compete to provide the best model to fit the data. A simple project to get started with Kaggle is the Titanic data set. Most Kaggle competitions end on a specific date. Website organizers have scheduled the Titanic competition to end on December 31, 20xx (with the year usually rolling forward). However, they have already extended the deadline several times, and an extension beyond 2014 is also possible. Second, the Titanic data set is considered a tutorial data set. There is no prize, and your score in the competition does not count towards becoming a Kaggle Master.

#### Kaggle Ranks

You achieve Kaggle ranks by earning gold, silver, and bronze medals.

- Kaggle Top Users
- Current Top Kaggle User's Profile Page
- Jeff Heaton's (your instructor) Kaggle Profile

#### Typical Kaggle Competition

A typical Kaggle competition will have several components. Consider the Titanic tutorial:

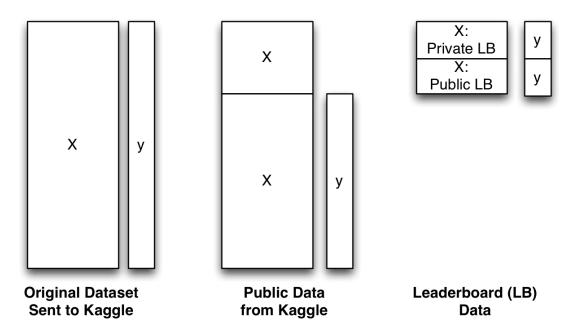
- Competition Summary Page
- Data Page
- Evaluation Description Page
- Leaderboard

#### How Kaggle Competition Scoring

Kaggle is provided with a data set by the competition sponsor, as seen in Figure 8.SCORE. Kaggle divides this data set as follows:

- Complete Data Set This is the complete data set.
  - Training Data Set This dataset provides both the inputs and the outcomes for the training portion of the data set.
  - Test Data Set This dataset provides the complete test data; however, it does not give the outcomes. Your submission file should contain the predicted results for this data set.
    - Public Leaderboard Kaggle does not tell you what part of the test data set contributes to the public leaderboard. Your public score is calculated based on this part of the data set.
    - Private Leaderboard Likewise, Kaggle does not tell you what part
      of the test data set contributes to the public leaderboard. Your final
      score/rank is calculated based on this part. You do not see your
      private leaderboard score until the end.

Figure 8.SCORE: How Kaggle Competition Scoring



## Preparing a Kaggle Submission

You do not submit the code to your solution to Kaggle. For competitions, you are scored entirely on the accuracy of your submission file. A Kaggle submission file is always a CSV file that contains the **Id** of the row you are predicting and the answer. For the titanic competition, a submission file looks something like this:

PassengerId, Survived 892,0 893,1 894,1 895,0 896,0 897,1

The above file states the prediction for each of the various passengers. You should only predict on ID's that are in the test file. Likewise, you should render a prediction for every row in the test file. Some competitions will have different formats for their answers. For example, a multi-classification will usually have a column for each class and your predictions for each class.

#### Select Kaggle Competitions

There have been many exciting competitions on Kaggle; these are some of my favorites. Some select predictive modeling competitions which use tabular data include:

- Otto Group Product Classification Challenge
- Galaxy Zoo The Galaxy Challenge
- Practice Fusion Diabetes Classification
- Predicting a Biological Response

Many Kaggle competitions include computer vision datasets, such as:

- Diabetic Retinopathy Detection
- Cats vs Dogs
- State Farm Distracted Driver Detection

## Module 8 Assignment

You can find the first assignment here: assignment 8

In [ ]:



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## Google CoLab Instructions

The following code ensures that Google CoLab is running the correct version of TensorFlow. Running the following code will map your GDrive to /content/drive.

```
s = sec_elapsed % 60
return "{}:{:>02}:{:>05.2f}".format(h, m, s)
```

Mounted at /content/drive Note: using Google CoLab

# Part 8.2: Building Ensembles with Scikit-Learn and Keras

#### **Evaluating Feature Importance**

Feature importance tells us how important each feature (from the feature/import vector) is to predicting a neural network or another model. There are many different ways to evaluate the feature importance of neural networks. The following paper presents an excellent (and readable) overview of the various means of assessing the significance of neural network inputs/features.

 An accurate comparison of methods for quantifying variable importance in artificial neural networks using simulated data [Cite:olden2004accurate].
 Ecological Modelling, 178(3), 389-397.

In summary, the following methods are available to neural networks:

- Connection Weights Algorithm
- Partial Derivatives
- Input Perturbation
- Sensitivity Analysis
- Forward Stepwise Addition
- Improved Stepwise Selection 1
- Backward Stepwise Elimination
- Improved Stepwise Selection

For this chapter, we will use the input Perturbation feature ranking algorithm. This algorithm will work with any regression or classification network. In the next section, I provide an implementation of the input perturbation algorithm for scikit-learn. This code implements a function below that will work with any scikit-learn model.

Leo Breiman provided this algorithm in his seminal paper on random forests. [Citebreiman2001random:] Although he presented this algorithm in conjunction with random forests, it is model-independent and appropriate for any supervised learning model. This algorithm, known as the input perturbation algorithm, works by evaluating a trained model's accuracy with each input individually shuffled from a data set. Shuffling an input causes it to become useless—effectively removing it from the model. More important inputs will produce a less accurate

score when they are removed by shuffling them. This process makes sense because important features will contribute to the model's accuracy. I first presented the TensorFlow implementation of this algorithm in the following paper.

 Early stabilizing feature importance for TensorFlow deep neural networks[Cite:heaton2017early]

This algorithm will use log loss to evaluate a classification problem and RMSE for regression.

```
In [2]: from sklearn import metrics
        import scipy as sp
        import numpy as np
        import math
        from sklearn import metrics
        def perturbation_rank(model, x, y, names, regression):
            errors = []
            for i in range(x.shape[1]):
                hold = np.array(x[:, i])
                np.random.shuffle(x[:, i])
                if regression:
                   pred = model.predict(x)
                    error = metrics.mean squared error(y, pred)
                else:
                    pred = model.predict(x)
                    error = metrics.log_loss(y, pred)
                errors.append(error)
                x[:, i] = hold
            max_error = np.max(errors)
            importance = [e/max error for e in errors]
            data = {'name':names,'error':errors,'importance':importance}
            result = pd.DataFrame(data, columns = ['name','error','importance'])
            result.sort values(by=['importance'], ascending=[0], inplace=True)
            result.reset index(inplace=True, drop=True)
            return result
```

### Classification and Input Perturbation Ranking

We now look at the code to perform perturbation ranking for a classification neural network. The implementation technique is slightly different for classification vs. regression, so I must provide two different implementations. The primary difference between classification and regression is how we evaluate the accuracy of the neural network in each of these two network types. We will

use the Root Mean Square (RMSE) error calculation, whereas we will use log loss for classification.

The code presented below creates a classification neural network that will predict the classic iris dataset.

```
In [3]: # HIDE OUTPUT
        import pandas as pd
        import io
        import requests
        import numpy as np
        from sklearn import metrics
        from tensorflow.keras.models import Sequential
        from tensorflow.keras.layers import Dense, Activation
        from tensorflow.keras.callbacks import EarlyStopping
        from sklearn.model selection import train test split
        df = pd.read csv(
            "https://data.heatonresearch.com/data/t81-558/iris.csv",
            na values=['NA', '?'])
        # Convert to numpy - Classification
        x = df[['sepal l', 'sepal w', 'petal l', 'petal w']].values
        dummies = pd.get dummies(df['species']) # Classification
        species = dummies.columns
        y = dummies.values
        # Split into train/test
        x train, x test, y train, y test = train test split(
            x, y, test size=0.25, random state=42)
        # Build neural network
        model = Sequential()
        model.add(Dense(50, input dim=x.shape[1], activation='relu')) # Hidden 1
        model.add(Dense(25, activation='relu')) # Hidden 2
        model.add(Dense(y.shape[1],activation='softmax')) # Output
        model.compile(loss='categorical crossentropy', optimizer='adam')
        model.fit(x train,y train,verbose=2,epochs=100)
```

```
Epoch 1/100
4/4 - 1s - loss: 2.0814 - 1s/epoch - 292ms/step
Epoch 2/100
4/4 - 0s - loss: 1.6125 - 14ms/epoch - 4ms/step
Epoch 3/100
4/4 - 0s - loss: 1.3316 - 26ms/epoch - 7ms/step
Epoch 4/100
4/4 - 0s - loss: 1.2246 - 13ms/epoch - 3ms/step
Epoch 5/100
4/4 - 0s - loss: 1.1989 - 13ms/epoch - 3ms/step
Epoch 6/100
4/4 - 0s - loss: 1.1349 - 14ms/epoch - 4ms/step
Epoch 7/100
4/4 - 0s - loss: 1.0543 - 21ms/epoch - 5ms/step
Epoch 8/100
4/4 - 0s - loss: 0.9987 - 25ms/epoch - 6ms/step
Epoch 9/100
4/4 - 0s - loss: 0.9449 - 20ms/epoch - 5ms/step
Epoch 10/100
4/4 - 0s - loss: 0.9032 - 16ms/epoch - 4ms/step
Epoch 11/100
4/4 - 0s - loss: 0.8623 - 20ms/epoch - 5ms/step
Epoch 12/100
4/4 - 0s - loss: 0.8274 - 12ms/epoch - 3ms/step
Epoch 13/100
4/4 - 0s - loss: 0.8013 - 18ms/epoch - 4ms/step
Epoch 14/100
4/4 - 0s - loss: 0.7718 - 18ms/epoch - 5ms/step
Epoch 15/100
4/4 - 0s - loss: 0.7426 - 19ms/epoch - 5ms/step
Epoch 16/100
4/4 - 0s - loss: 0.7163 - 13ms/epoch - 3ms/step
Epoch 17/100
4/4 - 0s - loss: 0.6933 - 13ms/epoch - 3ms/step
Epoch 18/100
4/4 - 0s - loss: 0.6689 - 14ms/epoch - 3ms/step
Epoch 19/100
4/4 - 0s - loss: 0.6488 - 11ms/epoch - 3ms/step
Epoch 20/100
4/4 - 0s - loss: 0.6294 - 11ms/epoch - 3ms/step
Epoch 21/100
4/4 - 0s - loss: 0.6094 - 20ms/epoch - 5ms/step
Epoch 22/100
4/4 - 0s - loss: 0.5911 - 18ms/epoch - 4ms/step
Epoch 23/100
4/4 - 0s - loss: 0.5725 - 16ms/epoch - 4ms/step
Epoch 24/100
4/4 - 0s - loss: 0.5550 - 13ms/epoch - 3ms/step
Epoch 25/100
4/4 - 0s - loss: 0.5389 - 14ms/epoch - 3ms/step
Epoch 26/100
4/4 - 0s - loss: 0.5207 - 15ms/epoch - 4ms/step
Epoch 27/100
4/4 - 0s - loss: 0.5041 - 14ms/epoch - 4ms/step
Epoch 28/100
4/4 - 0s - loss: 0.4901 - 14ms/epoch - 3ms/step
```

```
Epoch 29/100
4/4 - 0s - loss: 0.4765 - 14ms/epoch - 4ms/step
Epoch 30/100
4/4 - 0s - loss: 0.4619 - 16ms/epoch - 4ms/step
Epoch 31/100
4/4 - 0s - loss: 0.4489 - 16ms/epoch - 4ms/step
Epoch 32/100
4/4 - 0s - loss: 0.4366 - 13ms/epoch - 3ms/step
Epoch 33/100
4/4 - 0s - loss: 0.4243 - 13ms/epoch - 3ms/step
Epoch 34/100
4/4 - 0s - loss: 0.4124 - 14ms/epoch - 3ms/step
Epoch 35/100
4/4 - 0s - loss: 0.4015 - 14ms/epoch - 3ms/step
Epoch 36/100
4/4 - 0s - loss: 0.3917 - 21ms/epoch - 5ms/step
Epoch 37/100
4/4 - 0s - loss: 0.3826 - 30ms/epoch - 7ms/step
Epoch 38/100
4/4 - 0s - loss: 0.3713 - 18ms/epoch - 4ms/step
Epoch 39/100
4/4 - 0s - loss: 0.3621 - 16ms/epoch - 4ms/step
Epoch 40/100
4/4 - 0s - loss: 0.3543 - 14ms/epoch - 4ms/step
Epoch 41/100
4/4 - 0s - loss: 0.3460 - 14ms/epoch - 4ms/step
Epoch 42/100
4/4 - 0s - loss: 0.3385 - 28ms/epoch - 7ms/step
Epoch 43/100
4/4 - 0s - loss: 0.3280 - 31ms/epoch - 8ms/step
Epoch 44/100
4/4 - 0s - loss: 0.3211 - 15ms/epoch - 4ms/step
Epoch 45/100
4/4 - 0s - loss: 0.3144 - 14ms/epoch - 3ms/step
Epoch 46/100
4/4 - 0s - loss: 0.3068 - 15ms/epoch - 4ms/step
Epoch 47/100
4/4 - 0s - loss: 0.2992 - 19ms/epoch - 5ms/step
Epoch 48/100
4/4 - 0s - loss: 0.2922 - 18ms/epoch - 4ms/step
Epoch 49/100
4/4 - 0s - loss: 0.2847 - 37ms/epoch - 9ms/step
Epoch 50/100
4/4 - 0s - loss: 0.2803 - 13ms/epoch - 3ms/step
Epoch 51/100
4/4 - 0s - loss: 0.2756 - 13ms/epoch - 3ms/step
Epoch 52/100
4/4 - 0s - loss: 0.2665 - 18ms/epoch - 4ms/step
Epoch 53/100
4/4 - 0s - loss: 0.2632 - 16ms/epoch - 4ms/step
Epoch 54/100
4/4 - 0s - loss: 0.2571 - 20ms/epoch - 5ms/step
Epoch 55/100
4/4 - 0s - loss: 0.2499 - 17ms/epoch - 4ms/step
Epoch 56/100
4/4 - 0s - loss: 0.2458 - 17ms/epoch - 4ms/step
```

```
Epoch 57/100
4/4 - 0s - loss: 0.2399 - 23ms/epoch - 6ms/step
Epoch 58/100
4/4 - 0s - loss: 0.2340 - 16ms/epoch - 4ms/step
Epoch 59/100
4/4 - 0s - loss: 0.2318 - 16ms/epoch - 4ms/step
Epoch 60/100
4/4 - 0s - loss: 0.2225 - 12ms/epoch - 3ms/step
Epoch 61/100
4/4 - 0s - loss: 0.2266 - 15ms/epoch - 4ms/step
Epoch 62/100
4/4 - 0s - loss: 0.2178 - 12ms/epoch - 3ms/step
Epoch 63/100
4/4 - 0s - loss: 0.2116 - 15ms/epoch - 4ms/step
Epoch 64/100
4/4 - 0s - loss: 0.2137 - 21ms/epoch - 5ms/step
Epoch 65/100
4/4 - 0s - loss: 0.2030 - 17ms/epoch - 4ms/step
Epoch 66/100
4/4 - 0s - loss: 0.2041 - 14ms/epoch - 3ms/step
Epoch 67/100
4/4 - 0s - loss: 0.2001 - 15ms/epoch - 4ms/step
Epoch 68/100
4/4 - 0s - loss: 0.1919 - 25ms/epoch - 6ms/step
Epoch 69/100
4/4 - 0s - loss: 0.1894 - 23ms/epoch - 6ms/step
Epoch 70/100
4/4 - 0s - loss: 0.1863 - 17ms/epoch - 4ms/step
Epoch 71/100
4/4 - 0s - loss: 0.1823 - 16ms/epoch - 4ms/step
Epoch 72/100
4/4 - 0s - loss: 0.1790 - 24ms/epoch - 6ms/step
Epoch 73/100
4/4 - 0s - loss: 0.1780 - 16ms/epoch - 4ms/step
Epoch 74/100
4/4 - 0s - loss: 0.1755 - 15ms/epoch - 4ms/step
Epoch 75/100
4/4 - 0s - loss: 0.1719 - 31ms/epoch - 8ms/step
Epoch 76/100
4/4 - 0s - loss: 0.1767 - 21ms/epoch - 5ms/step
Epoch 77/100
4/4 - 0s - loss: 0.1694 - 17ms/epoch - 4ms/step
Epoch 78/100
4/4 - 0s - loss: 0.1655 - 27ms/epoch - 7ms/step
Epoch 79/100
4/4 - 0s - loss: 0.1634 - 23ms/epoch - 6ms/step
Epoch 80/100
4/4 - 0s - loss: 0.1566 - 17ms/epoch - 4ms/step
Epoch 81/100
4/4 - 0s - loss: 0.1563 - 17ms/epoch - 4ms/step
Epoch 82/100
4/4 - 0s - loss: 0.1536 - 15ms/epoch - 4ms/step
Epoch 83/100
4/4 - 0s - loss: 0.1504 - 18ms/epoch - 4ms/step
Epoch 84/100
4/4 - 0s - loss: 0.1502 - 16ms/epoch - 4ms/step
```

```
Epoch 85/100
       4/4 - 0s - loss: 0.1469 - 17ms/epoch - 4ms/step
       Epoch 86/100
       4/4 - 0s - loss: 0.1448 - 28ms/epoch - 7ms/step
       Epoch 87/100
       4/4 - 0s - loss: 0.1424 - 23ms/epoch - 6ms/step
       Epoch 88/100
       4/4 - 0s - loss: 0.1401 - 25ms/epoch - 6ms/step
       Epoch 89/100
       4/4 - 0s - loss: 0.1386 - 47ms/epoch - 12ms/step
       Epoch 90/100
       4/4 - 0s - loss: 0.1365 - 30ms/epoch - 7ms/step
       Epoch 91/100
       4/4 - 0s - loss: 0.1383 - 41ms/epoch - 10ms/step
       Epoch 92/100
       4/4 - 0s - loss: 0.1332 - 12ms/epoch - 3ms/step
       Epoch 93/100
       4/4 - 0s - loss: 0.1311 - 20ms/epoch - 5ms/step
       Epoch 94/100
       4/4 - 0s - loss: 0.1320 - 20ms/epoch - 5ms/step
       Epoch 95/100
       4/4 - 0s - loss: 0.1302 - 17ms/epoch - 4ms/step
       Epoch 96/100
       4/4 - 0s - loss: 0.1311 - 19ms/epoch - 5ms/step
       Epoch 97/100
       4/4 - 0s - loss: 0.1248 - 14ms/epoch - 3ms/step
       Epoch 98/100
       4/4 - 0s - loss: 0.1254 - 12ms/epoch - 3ms/step
       Epoch 99/100
       4/4 - 0s - loss: 0.1275 - 13ms/epoch - 3ms/step
       Epoch 100/100
       4/4 - 0s - loss: 0.1225 - 41ms/epoch - 10ms/step
Out[3]: <keras.callbacks.History at 0x7fc869fd2950>
```

Next, we evaluate the accuracy of the trained model. Here we see that the neural network performs great, with an accuracy of 1.0. We might fear overfitting with such high accuracy for a more complex dataset. However, for this example, we are more interested in determining the importance of each column.

```
In [4]: from sklearn.metrics import accuracy_score

pred = model.predict(x_test)
predict_classes = np.argmax(pred,axis=1)
expected_classes = np.argmax(y_test,axis=1)
correct = accuracy_score(expected_classes,predict_classes)
print(f"Accuracy: {correct}")
```

Accuracy: 1.0

We are now ready to call the input perturbation algorithm. First, we extract the column names and remove the target column. The target column is not important, as it is the objective, not one of the inputs. In supervised learning, the target is of the utmost importance.

We can see the importance displayed in the following table. The most important column is always 1.0, and lessor columns will continue in a downward trend. The least important column will have the lowest rank.

```
In [5]: # Rank the features
    from IPython.display import display, HTML

names = list(df.columns) # x+y column names
names.remove("species") # remove the target(y)
rank = perturbation_rank(model, x_test, y_test, names, False)
display(rank)
```

	name	error	importance
0	petal_l	2.609378	1.000000
1	petal_w	0.480387	0.184100
2	sepal_l	0.223239	0.085553
3	sepal_w	0.128518	0.049252

### Regression and Input Perturbation Ranking

We now see how to use input perturbation ranking for a regression neural network. We will use the MPG dataset as a demonstration. The code below loads the MPG dataset and creates a regression neural network for this dataset. The code trains the neural network and calculates an RMSE evaluation.

```
In [6]: # HIDE OUTPUT
        from tensorflow.keras.models import Sequential
        from tensorflow.keras.layers import Dense, Activation
        from sklearn.model selection import train test split
        import pandas as pd
        import io
        import os
        import requests
        import numpy as np
        from sklearn import metrics
        save path = "."
        df = pd.read csv(
            "https://data.heatonresearch.com/data/t81-558/auto-mpg.csv",
            na values=['NA', '?'])
        cars = df['name']
        # Handle missing value
        df['horsepower'] = df['horsepower'].fillna(df['horsepower'].median())
        # Pandas to Numpy
```

```
Epoch 1/100
10/10 - 1s - loss: 328433.8125 - 898ms/epoch - 90ms/step
Epoch 2/100
10/10 - 0s - loss: 78914.6406 - 26ms/epoch - 3ms/step
Epoch 3/100
10/10 - 0s - loss: 5371.1025 - 50ms/epoch - 5ms/step
Epoch 4/100
10/10 - 0s - loss: 4021.9753 - 34ms/epoch - 3ms/step
Epoch 5/100
10/10 - 0s - loss: 4438.5728 - 33ms/epoch - 3ms/step
Epoch 6/100
10/10 - 0s - loss: 1030.3115 - 34ms/epoch - 3ms/step
Epoch 7/100
10/10 - 0s - loss: 594.9177 - 31ms/epoch - 3ms/step
Epoch 8/100
10/10 - 0s - loss: 655.3908 - 31ms/epoch - 3ms/step
Epoch 9/100
10/10 - 0s - loss: 465.0457 - 25ms/epoch - 2ms/step
Epoch 10/100
10/10 - 0s - loss: 458.7520 - 30ms/epoch - 3ms/step
Epoch 11/100
10/10 - 0s - loss: 452.4102 - 22ms/epoch - 2ms/step
Epoch 12/100
10/10 - 0s - loss: 439.8730 - 25ms/epoch - 3ms/step
Epoch 13/100
10/10 - 0s - loss: 434.8245 - 27ms/epoch - 3ms/step
Epoch 14/100
10/10 - 0s - loss: 433.7303 - 25ms/epoch - 3ms/step
Epoch 15/100
10/10 - 0s - loss: 427.2859 - 46ms/epoch - 5ms/step
Epoch 16/100
10/10 - 0s - loss: 424.1164 - 50ms/epoch - 5ms/step
Epoch 17/100
10/10 - 0s - loss: 422.3007 - 42ms/epoch - 4ms/step
Epoch 18/100
10/10 - 0s - loss: 418.4877 - 31ms/epoch - 3ms/step
Epoch 19/100
10/10 - 0s - loss: 414.2283 - 23ms/epoch - 2ms/step
Epoch 20/100
10/10 - 0s - loss: 410.2691 - 34ms/epoch - 3ms/step
Epoch 21/100
10/10 - 0s - loss: 407.0490 - 29ms/epoch - 3ms/step
Epoch 22/100
10/10 - 0s - loss: 406.2433 - 46ms/epoch - 5ms/step
Epoch 23/100
10/10 - 0s - loss: 399.7404 - 37ms/epoch - 4ms/step
Epoch 24/100
10/10 - 0s - loss: 396.3280 - 66ms/epoch - 7ms/step
Epoch 25/100
10/10 - 0s - loss: 391.0629 - 28ms/epoch - 3ms/step
Epoch 26/100
10/10 - 0s - loss: 387.3203 - 26ms/epoch - 3ms/step
Epoch 27/100
10/10 - 0s - loss: 382.7670 - 54ms/epoch - 5ms/step
Epoch 28/100
10/10 - 0s - loss: 380.6316 - 21ms/epoch - 2ms/step
```

```
Epoch 29/100
10/10 - 0s - loss: 375.9518 - 30ms/epoch - 3ms/step
Epoch 30/100
10/10 - 0s - loss: 372.7001 - 24ms/epoch - 2ms/step
Epoch 31/100
10/10 - 0s - loss: 366.7871 - 24ms/epoch - 2ms/step
Epoch 32/100
10/10 - 0s - loss: 363.4180 - 42ms/epoch - 4ms/step
Epoch 33/100
10/10 - 0s - loss: 359.6006 - 47ms/epoch - 5ms/step
Epoch 34/100
10/10 - 0s - loss: 359.4055 - 46ms/epoch - 5ms/step
Epoch 35/100
10/10 - 0s - loss: 350.7181 - 29ms/epoch - 3ms/step
Epoch 36/100
10/10 - 0s - loss: 348.6260 - 42ms/epoch - 4ms/step
Epoch 37/100
10/10 - 0s - loss: 343.6122 - 28ms/epoch - 3ms/step
Epoch 38/100
10/10 - 0s - loss: 339.6165 - 32ms/epoch - 3ms/step
Epoch 39/100
10/10 - 0s - loss: 334.5634 - 32ms/epoch - 3ms/step
Epoch 40/100
10/10 - 0s - loss: 332.6061 - 34ms/epoch - 3ms/step
Epoch 41/100
10/10 - 0s - loss: 326.7434 - 22ms/epoch - 2ms/step
Epoch 42/100
10/10 - 0s - loss: 323.8063 - 40ms/epoch - 4ms/step
Epoch 43/100
10/10 - 0s - loss: 320.2585 - 29ms/epoch - 3ms/step
Epoch 44/100
10/10 - 0s - loss: 315.3609 - 23ms/epoch - 2ms/step
Epoch 45/100
10/10 - 0s - loss: 311.4920 - 23ms/epoch - 2ms/step
Epoch 46/100
10/10 - 0s - loss: 308.9212 - 29ms/epoch - 3ms/step
Epoch 47/100
10/10 - 0s - loss: 303.1410 - 24ms/epoch - 2ms/step
Epoch 48/100
10/10 - 0s - loss: 299.9317 - 24ms/epoch - 2ms/step
Epoch 49/100
10/10 - 0s - loss: 294.4305 - 23ms/epoch - 2ms/step
Epoch 50/100
10/10 - 0s - loss: 291.4469 - 24ms/epoch - 2ms/step
Epoch 51/100
10/10 - 0s - loss: 287.3263 - 41ms/epoch - 4ms/step
Epoch 52/100
10/10 - 0s - loss: 284.3096 - 49ms/epoch - 5ms/step
Epoch 53/100
10/10 - 0s - loss: 280.5522 - 30ms/epoch - 3ms/step
Epoch 54/100
10/10 - 0s - loss: 276.1487 - 26ms/epoch - 3ms/step
Epoch 55/100
10/10 - 0s - loss: 271.3444 - 42ms/epoch - 4ms/step
Epoch 56/100
10/10 - 0s - loss: 280.0936 - 33ms/epoch - 3ms/step
```

```
Epoch 57/100
10/10 - 0s - loss: 263.7166 - 40ms/epoch - 4ms/step
Epoch 58/100
10/10 - 0s - loss: 261.6750 - 56ms/epoch - 6ms/step
Epoch 59/100
10/10 - 0s - loss: 258.5714 - 45ms/epoch - 4ms/step
Epoch 60/100
10/10 - 0s - loss: 252.6791 - 31ms/epoch - 3ms/step
Epoch 61/100
10/10 - 0s - loss: 250.1348 - 53ms/epoch - 5ms/step
Epoch 62/100
10/10 - 0s - loss: 246.4157 - 72ms/epoch - 7ms/step
Epoch 63/100
10/10 - 0s - loss: 242.3768 - 46ms/epoch - 5ms/step
Epoch 64/100
10/10 - 0s - loss: 238.7874 - 28ms/epoch - 3ms/step
Epoch 65/100
10/10 - 0s - loss: 235.8578 - 42ms/epoch - 4ms/step
Epoch 66/100
10/10 - 0s - loss: 233.7492 - 24ms/epoch - 2ms/step
Epoch 67/100
10/10 - 0s - loss: 229.0066 - 26ms/epoch - 3ms/step
Epoch 68/100
10/10 - 0s - loss: 225.7449 - 25ms/epoch - 3ms/step
Epoch 69/100
10/10 - 0s - loss: 223.5038 - 25ms/epoch - 2ms/step
Epoch 70/100
10/10 - 0s - loss: 219.9561 - 39ms/epoch - 4ms/step
Epoch 71/100
10/10 - 0s - loss: 215.1055 - 58ms/epoch - 6ms/step
Epoch 72/100
10/10 - 0s - loss: 211.9364 - 39ms/epoch - 4ms/step
Epoch 73/100
10/10 - 0s - loss: 208.1019 - 55ms/epoch - 5ms/step
Epoch 74/100
10/10 - 0s - loss: 207.4119 - 34ms/epoch - 3ms/step
Epoch 75/100
10/10 - 0s - loss: 206.8693 - 40ms/epoch - 4ms/step
Epoch 76/100
10/10 - 0s - loss: 197.9749 - 49ms/epoch - 5ms/step
Epoch 77/100
10/10 - 0s - loss: 196.9090 - 34ms/epoch - 3ms/step
Epoch 78/100
10/10 - 0s - loss: 192.6349 - 45ms/epoch - 4ms/step
Epoch 79/100
10/10 - 0s - loss: 189.6783 - 31ms/epoch - 3ms/step
Epoch 80/100
10/10 - 0s - loss: 186.6584 - 25ms/epoch - 2ms/step
Epoch 81/100
10/10 - 0s - loss: 186.1920 - 29ms/epoch - 3ms/step
Epoch 82/100
10/10 - 0s - loss: 181.1735 - 31ms/epoch - 3ms/step
Epoch 83/100
10/10 - 0s - loss: 177.9338 - 51ms/epoch - 5ms/step
Epoch 84/100
10/10 - 0s - loss: 174.6662 - 87ms/epoch - 9ms/step
```

```
Epoch 85/100
10/10 - 0s - loss: 172.9421 - 90ms/epoch - 9ms/step
Epoch 86/100
10/10 - 0s - loss: 169.1906 - 58ms/epoch - 6ms/step
Epoch 87/100
10/10 - 0s - loss: 166.4181 - 57ms/epoch - 6ms/step
Epoch 88/100
10/10 - 0s - loss: 163.7466 - 36ms/epoch - 4ms/step
Epoch 89/100
10/10 - 0s - loss: 161.4653 - 29ms/epoch - 3ms/step
Epoch 90/100
10/10 - 0s - loss: 158.6274 - 30ms/epoch - 3ms/step
Epoch 91/100
10/10 - 0s - loss: 159.4237 - 32ms/epoch - 3ms/step
Epoch 92/100
10/10 - 0s - loss: 159.2035 - 31ms/epoch - 3ms/step
Epoch 93/100
10/10 - 0s - loss: 150.2793 - 38ms/epoch - 4ms/step
Epoch 94/100
10/10 - 0s - loss: 148.9276 - 36ms/epoch - 4ms/step
Epoch 95/100
10/10 - 0s - loss: 146.7706 - 34ms/epoch - 3ms/step
Epoch 96/100
10/10 - 0s - loss: 144.4946 - 29ms/epoch - 3ms/step
Epoch 97/100
10/10 - 0s - loss: 141.5782 - 28ms/epoch - 3ms/step
Epoch 98/100
10/10 - 0s - loss: 139.3355 - 27ms/epoch - 3ms/step
Epoch 99/100
10/10 - 0s - loss: 136.9762 - 56ms/epoch - 6ms/step
Epoch 100/100
10/10 - 0s - loss: 135.6660 - 23ms/epoch - 2ms/step
```

Just as before, we extract the column names and discard the target. We can now create a ranking of the importance of each of the input features. The feature with a ranking of 1.0 is the most important.

```
In [7]: # Rank the features
    from IPython.display import display, HTML

    names = list(df.columns) # x+y column names
    names.remove("name")
    names.remove("mpg") # remove the target(y)
    rank = perturbation_rank(model, x_test, y_test, names, True)
    display(rank)
```

	name	error	importance
0	displacement	139.657598	1.000000
1	acceleration	139.261508	0.997164
2	origin	134.637690	0.964056
3	year	134.177126	0.960758
4	cylinders	132.747246	0.950519
5	horsepower	121.501102	0.869993
6	weight	75.244610	0.538779

#### Biological Response with Neural Network

The following sections will demonstrate how to use feature importance ranking and ensembling with a more complex dataset. Ensembling is the process where you combine multiple models for greater accuracy. Kaggle competition winners frequently make use of ensembling for high-ranking solutions.

We will use the biological response dataset, a Kaggle dataset, where there is an unusually high number of columns. Because of the large number of columns, it is essential to use feature ranking to determine the importance of these columns. We begin by loading the dataset and preprocessing. This Kaggle dataset is a binary classification problem. You must predict if certain conditions will cause a biological response.

Predicting a Biological Response

```
In [8]: import pandas as pd
import os
import numpy as np
from sklearn import metrics
from scipy.stats import zscore
from sklearn.model_selection import KFold
from IPython.display import HTML, display

URL = "https://data.heatonresearch.com/data/t81-558/kaggle/"

df_train = pd.read_csv(
    URL+"bio_train.csv",
    na_values=['NA', '?'])

df_test = pd.read_csv(
    URL+"bio_test.csv",
    na_values=['NA', '?'])

activity_classes = df_train['Activity']
```

A large number of columns is evident when we display the shape of the dataset.

The following code constructs a classification neural network and trains it for the biological response dataset. Once trained, the accuracy is measured.

```
In [10]: import os
         import pandas as pd
         import tensorflow as tf
         from tensorflow.keras.models import Sequential
         from tensorflow.keras.layers import Dense, Activation
         from sklearn.model selection import train test split
         from tensorflow.keras.callbacks import EarlyStopping
         import numpy as np
         import sklearn
         # Encode feature vector
         # Convert to numpy - Classification
         x columns = df train.columns.drop('Activity')
         x = df train[x columns].values
         y = df train['Activity'].values # Classification
         x submit = df test[x columns].values.astype(np.float32)
         # Split into train/test
         x train, x test, y train, y test = train test split(
             x, y, test size=0.25, random state=42)
         print("Fitting/Training...")
         model = Sequential()
         model.add(Dense(25, input dim=x.shape[1], activation='relu'))
         model.add(Dense(10))
         model.add(Dense(1,activation='sigmoid'))
         model.compile(loss='binary_crossentropy', optimizer='adam')
         monitor = EarlyStopping(monitor='val loss', min delta=1e-3,
                                 patience=5, verbose=1, mode='auto')
         model.fit(x_train,y_train,validation_data=(x_test,y_test),
                   callbacks=[monitor], verbose=0, epochs=1000)
         print("Fitting done...")
         # Predict
         pred = model.predict(x test).flatten()
         # Clip so that min is never exactly 0, max never 1
         pred = np.clip(pred,a min=le-6,a max=(1-le-6))
         print("Validation logloss: {}".format(
             sklearn.metrics.log loss(y test,pred)))
         # Evaluate success using accuracy
         pred = pred>0.5 # If greater than 0.5 probability, then true
         score = metrics.accuracy score(y test, pred)
```

#### What Features/Columns are Important

The following uses perturbation ranking to evaluate the neural network.

```
In [11]: # Rank the features
    from IPython.display import display, HTML

names = list(df_train.columns) # x+y column names
names.remove("Activity") # remove the target(y)
rank = perturbation_rank(model, x_test, y_test, names, False)
display(rank[0:10])
```

	name	error	importance
0	D27	0.603974	1.000000
1	D1049	0.565997	0.937122
2	D51	0.565883	0.936934
3	D998	0.563872	0.933604
4	D1059	0.563745	0.933394
5	D961	0.563723	0.933357
6	D1407	0.563532	0.933041
7	D1309	0.562244	0.930908
8	D1100	0.561902	0.930341
9	D1275	0.561659	0.929940

#### Neural Network Ensemble

A neural network ensemble combines neural network predictions with other models. The program determines the exact blend of these models by logistic regression. The following code performs this blend for a classification. If you present the final predictions from the ensemble to Kaggle, you will see that the result is very accurate.

```
In [12]: # HIDE OUTPUT
         import numpy as np
         import os
         import pandas as pd
         import math
         from tensorflow.keras.wrappers.scikit learn import KerasClassifier
         from sklearn.neighbors import KNeighborsClassifier
         from sklearn.model selection import StratifiedKFold
         from sklearn.ensemble import RandomForestClassifier
         from sklearn.ensemble import ExtraTreesClassifier
         from sklearn.ensemble import GradientBoostingClassifier
         from sklearn.linear model import LogisticRegression
         SHUFFLE = False
         FOLDS = 10
         def build ann(input size, classes, neurons):
             model = Sequential()
             model.add(Dense(neurons, input dim=input size, activation='relu'))
             model.add(Dense(1))
             model.add(Dense(classes,activation='softmax'))
             model.compile(loss='categorical crossentropy', optimizer='adam')
             return model
         def mlogloss(y test, preds):
             epsilon = 1e-15
             sum = 0
             for row in zip(preds,y_test):
                 x = row[0][row[1]]
                 x = max(epsilon, x)
                 x = min(1-epsilon,x)
                 sum+=math.log(x)
             return( (-1/len(preds))*sum)
         def stretch(y):
             return (y - y.min()) / (y.max() - y.min())
         def blend ensemble(x, y, x submit):
             kf = StratifiedKFold(FOLDS)
             folds = list(kf.split(x,y))
             models = [
                 KerasClassifier(build fn=build ann,neurons=20,
                             input size=x.shape[1],classes=2),
                 KNeighborsClassifier(n neighbors=3),
                 RandomForestClassifier(n estimators=100, n jobs=-1,
                                         criterion='gini'),
```

```
RandomForestClassifier(n estimators=100, n jobs=-1,
                               criterion='entropy'),
        ExtraTreesClassifier(n estimators=100, n jobs=-1,
                             criterion='gini'),
        ExtraTreesClassifier(n_estimators=100, n_jobs=-1,
                             criterion='entropy'),
        GradientBoostingClassifier(learning rate=0.05,
                subsample=0.5, max depth=6, n estimators=50)]
   dataset blend train = np.zeros((x.shape[0], len(models)))
   dataset blend test = np.zeros((x submit.shape[0], len(models)))
   for j, model in enumerate(models):
        print("Model: {} : {}".format(j, model) )
        fold sums = np.zeros((x submit.shape[0], len(folds)))
        total loss = 0
        for i, (train, test) in enumerate(folds):
            x train = x[train]
            y train = y[train]
            x \text{ test} = x[\text{test}]
            y \text{ test} = y[\text{test}]
            model.fit(x train, y train)
            pred = np.array(model.predict proba(x test))
            dataset blend train[test, j] = pred[:, 1]
            pred2 = np.array(model.predict proba(x submit))
            fold sums[:, i] = pred2[:, 1]
            loss = mlogloss(y test, pred)
            total loss+=loss
            print("Fold #{}: loss={}".format(i,loss))
        print("{}: Mean loss={}".format(model.__class__.__name__,
                                         total loss/len(folds)))
        dataset_blend_test[:, j] = fold_sums.mean(1)
   print()
   print("Blending models.")
   blend = LogisticRegression(solver='lbfgs')
   blend.fit(dataset blend train, y)
    return blend.predict proba(dataset blend test)
if name == ' main ':
   np.random.seed(42) # seed to shuffle the train set
   print("Loading data...")
   URL = "https://data.heatonresearch.com/data/t81-558/kaggle/"
   df train = pd.read csv(
        URL+"bio train.csv",
        na values=['NA', '?'])
   df submit = pd.read csv(
        URL+"bio test.csv",
        na values=['NA', '?'])
    predictors = list(df train.columns.values)
   predictors.remove('Activity')
```

```
x = df_train[predictors].values
y = df train['Activity']
x submit = df submit.values
if SHUFFLE:
    idx = np.random.permutation(y.size)
    x = x[idx]
    y = y[idx]
submit data = blend ensemble(x, y, x submit)
submit_data = stretch(submit_data)
###################
# Build submit file
#####################
ids = [id+1 for id in range(submit data.shape[0])]
submit df = pd.DataFrame({'MoleculeId': ids,
                          'PredictedProbability':
                          submit data[:, 1]},
                         columns=['MoleculeId',
                        'PredictedProbability'])
submit df.to csv("submit.csv", index=False)
```

#### Loading data...

/usr/local/lib/python3.7/dist-packages/ipykernel\_launcher.py:44: Deprecation Warning: KerasClassifier is deprecated, use Sci-Keras (https://github.com/adriangb/scikeras) instead. See https://www.adriangb.com/scikeras/stable/migration.html for help migrating.

```
Model: 0 : <keras.wrappers.scikit learn.KerasClassifier object at 0x7fc86980
9610>
Fold #0: loss=0.5544745638322883
Fold #1: loss=0.5684765604955473
Fold #2: loss=0.5214491621944897
106/106 [============= ] - 1s 2ms/step - loss: 0.6301
Fold #3: loss=0.5264746750391351
106/106 [============== ] - 1s 2ms/step - loss: 0.5905
Fold #4: loss=0.5327822461352748
106/106 [============] - 1s 2ms/step - loss: 0.5993
Fold #5: loss=0.5800157462831582
Fold #6: loss=0.5189563830365144
Fold #7: loss=0.5625417655617023
106/106 [=============] - 1s 2ms/step - loss: 0.5935
Fold #8: loss=0.5238374326475557
Fold #9: loss=0.5322226787930878
KerasClassifier: Mean loss=0.5421231214018752
Model: 1 : KNeighborsClassifier(n neighbors=3)
Fold #0: loss=3.606678388314123
Fold #1: loss=2.2256421551487593
Fold #2: loss=3.6815437059542186
Fold #3: loss=2.416161292225968
Fold #4: loss=4.442472310149748
Fold #5: loss=4.321350530738247
Fold #6: loss=3.400455469543658
Fold #7: loss=3.1724147110842513
Fold #8: loss=2.117356283193681
Fold #9: loss=3.0532135963322586
KNeighborsClassifier: Mean loss=3.243728844268491
Model: 2 : RandomForestClassifier(n jobs=-1)
Fold #0: loss=0.4657177982691548
Fold #1: loss=0.4346825805694879
Fold #2: loss=0.4593868993445528
Fold #3: loss=0.41674899522216713
Fold #4: loss=0.4851849131056564
Fold #5: loss=0.48473291073937
Fold #6: loss=0.41274608628217674
Fold #7: loss=0.47405291219252377
Fold #8: loss=0.44974230059938286
Fold #9: loss=0.46340159258241087
RandomForestClassifier: Mean loss=0.45463969889068834
Model: 3 : RandomForestClassifier(criterion='entropy', n jobs=-1)
Fold #0: loss=0.4511847247326708
Fold #1: loss=0.42707704254926593
Fold #2: loss=0.5550335199035183
Fold #3: loss=0.42186970733328516
Fold #4: loss=0.4794331756190797
Fold #5: loss=0.4730559509802762
Fold #6: loss=0.41116235817215196
Fold #7: loss=0.46835919493314265
```

```
Fold #8: loss=0.4496144890690015
Fold #9: loss=0.4625902934553457
RandomForestClassifier: Mean loss=0.4599380456747738
Model: 4 : ExtraTreesClassifier(n jobs=-1)
Fold #0: loss=0.45496751079363495
Fold #1: loss=0.5013051157905043
Fold #2: loss=0.5886179891724027
Fold #3: loss=0.41646902160044674
Fold #4: loss=0.4957910697444236
Fold #5: loss=0.4773401028797005
Fold #6: loss=0.41935061504547827
Fold #7: loss=0.5757908399174205
Fold #8: loss=0.4585195863412778
Fold #9: loss=0.6210675972963805
ExtraTreesClassifier: Mean loss=0.500921944858167
Model: 5 : ExtraTreesClassifier(criterion='entropy', n jobs=-1)
Fold #0: loss=0.44825346440152214
Fold #1: loss=0.40764412171784686
Fold #2: loss=0.5819367378417363
Fold #3: loss=0.4140589874942631
Fold #4: loss=0.4923489720481471
Fold #5: loss=0.5744429921555051
Fold #6: loss=0.42334390524742155
Fold #7: loss=0.6409291880353659
Fold #8: loss=0.45627884947155956
Fold #9: loss=0.466653395317917
ExtraTreesClassifier: Mean loss=0.49058906137312847
Model: 6 : GradientBoostingClassifier(learning_rate=0.05, max_depth=6, n_est
imators=50,
                           subsample=0.5)
Fold #0: loss=0.4789324034433162
Fold #1: loss=0.4565636914381977
Fold #2: loss=0.47057741836357014
Fold #3: loss=0.4436328438944843
Fold #4: loss=0.4883293501002484
Fold #5: loss=0.4843521206311074
Fold #6: loss=0.4436043855503229
Fold #7: loss=0.45977393398397765
Fold #8: loss=0.46632256794136323
Fold #9: loss=0.4703354072414907
GradientBoostingClassifier: Mean loss=0.4662424122588078
Blending models.
```



# T81-558: Applications of Deep Neural Networks

#### Module 8: Kaggle Data Sets

- Instructor: Jeff Heaton, McKelvey School of Engineering, Washington University in St. Louis
- For more information visit the class website.

#### Module 8 Material

- Part 8.1: Introduction to Kaggle [Video] [Notebook]
- Part 8.2: Building Ensembles with Scikit-Learn and Keras [Video] [Notebook]
- Part 8.3: How Should you Architect Your Keras Neural Network: Hyperparameters [Video] [Notebook]
- Part 8.4: Bayesian Hyperparameter Optimization for Keras [Video] [Notebook]
- Part 8.5: Current Semester's Kaggle [Video] [Notebook]

## Google CoLab Instructions

The following code ensures that Google CoLab is running the correct version of TensorFlow.

# Part 8.3: Architecting Network: Hyperparameters

You have probably noticed several hyperparameters introduced previously in this course that you need to choose for your neural network. The number of layers, neuron counts per layer, layer types, and activation functions are all choices you must make to optimize your neural network. Some of the categories of hyperparameters for you to choose from coming from the following list:

- Number of Hidden Layers and Neuron Counts
- Activation Functions
- Advanced Activation Functions
- Regularization: L1, L2, Dropout
- Batch Normalization
- Training Parameters

The following sections will introduce each of these categories for Keras. While I will provide some general guidelines for hyperparameter selection, no two tasks are the same. You will benefit from experimentation with these values to determine what works best for your neural network. In the next part, we will see how machine learning can select some of these values independently.

#### Number of Hidden Layers and Neuron Counts

The structure of Keras layers is perhaps the hyperparameters that most become aware of first. How many layers should you have? How many neurons are on each layer? What activation function and layer type should you use? These are all questions that come up when designing a neural network. There are many different types of layer in Keras, listed here:

- Activation You can also add activation functions as layers. Using the
  activation layer is the same as specifying the activation function as part of a
  Dense (or other) layer type.
- **ActivityRegularization** Used to add L1/L2 regularization outside of a layer. You can specify L1 and L2 as part of a Dense (or other) layer type.
- **Dense** The original neural network layer type. In this layer type, every neuron connects to the next layer. The input vector is one-dimensional, and placing specific inputs next does not affect each other.
- Dropout Dropout consists of randomly setting a fraction rate of input units to 0 at each update during training time, which helps prevent overfitting.
   Dropout only occurs during training.

- Flatten Flattens the input to 1D and does not affect the batch size.
- **Input** A Keras tensor is a tensor object from the underlying back end (Theano, TensorFlow, or CNTK), which we augment with specific attributes to build a Keras by knowing the inputs and outputs of the model.
- Lambda Wraps arbitrary expression as a Layer object.
- **Masking** Masks a sequence using a mask value to skip timesteps.
- **Permute** Permutes the input dimensions according to a given pattern. Useful for tasks such as connecting RNNs and convolutional networks.
- **RepeatVector** Repeats the input n times.
- **Reshape** Similar to Numpy reshapes.
- **SpatialDropout1D** This version performs the same function as Dropout; however, it drops entire 1D feature maps instead of individual elements.
- **SpatialDropout2D** This version performs the same function as Dropout; however, it drops entire 2D feature maps instead of individual elements
- **SpatialDropout3D** This version performs the same function as Dropout; however, it drops entire 3D feature maps instead of individual elements.

There is always trial and error for choosing a good number of neurons and hidden layers. Generally, the number of neurons on each layer will be larger closer to the hidden layer and smaller towards the output layer. This configuration gives the neural network a somewhat triangular or trapezoid appearance.

#### **Activation Functions**

Activation functions are a choice that you must make for each layer. Generally, you can follow this guideline:

- Hidden Layers RELU
- Output Layer Softmax for classification, linear for regression.

Some of the common activation functions in Keras are listed here:

- **softmax** Used for multi-class classification. Ensures all output neurons behave as probabilities and sum to 1.0.
- **elu** Exponential linear unit. Exponential Linear Unit or its widely known name ELU is a function that tends to converge cost to zero faster and produce more accurate results. Can produce negative outputs.
- **selu** Scaled Exponential Linear Unit (SELU), essentially **elu** multiplied by a scaling constant.
- **softplus** Softplus activation function. log(exp(x) + 1) Introduced in 2001.
- **softsign** Softsign activation function. x/(abs(x)+1) Similar to tanh, but not widely used.

- **relu** Very popular neural network activation function. Used for hidden layers, cannot output negative values. No trainable parameters.
- **tanh** Classic neural network activation function, though often replaced by relu family on modern networks.
- **sigmoid** Classic neural network activation. Often used on output layer of a binary classifier.
- hard\_sigmoid Less computationally expensive variant of sigmoid.
- exponential Exponential (base e) activation function.
- **linear** Pass-through activation function. Usually used on the output layer of a regression neural network.

For more information about Keras activation functions refer to the following:

- Keras Activation Functions
- Activation Function Cheat Sheets

#### **Advanced Activation Functions**

Hyperparameters are not changed when the neural network trains. You, the network designer, must define the hyperparameters. The neural network learns regular parameters during neural network training. Neural network weights are the most common type of regular parameter. The "advanced activation functions," as Keras call them, also contain parameters that the network will learn during training. These activation functions may give you better performance than RELU.

- **LeakyReLU** Leaky version of a Rectified Linear Unit. It allows a small gradient when the unit is not active, controlled by alpha hyperparameter.
- **PReLU** Parametric Rectified Linear Unit, learns the alpha hyperparameter.

#### Regularization: L1, L2, Dropout

- Keras Regularization
- Keras Dropout

#### **Batch Normalization**

- Keras Batch Normalization
- Ioffe, S., & Szegedy, C. (2015). Batch normalization: Accelerating deep network training by reducing internal covariate shift. *arXiv* preprint *arXiv*:1502.03167.

Normalize the activations of the previous layer at each batch, i.e. applies a transformation that maintains the mean activation close to 0 and the activation standard deviation close to 1. Can allow learning rate to be larger.

## **Training Parameters**

- Keras Optimizers
- Batch Size Usually small, such as 32 or so.
- Learning Rate Usually small, 1e-3 or so.

In [ ]:



# T81-558: Applications of Deep Neural Networks

#### Module 8: Kaggle Data Sets

- Instructor: Jeff Heaton, McKelvey School of Engineering, Washington University in St. Louis
- For more information visit the class website.

#### Module 8 Material

- Part 8.1: Introduction to Kaggle [Video] [Notebook]
- Part 8.2: Building Ensembles with Scikit-Learn and Keras [Video] [Notebook]
- Part 8.3: How Should you Architect Your Keras Neural Network: Hyperparameters [Video] [Notebook]
- Part 8.4: Bayesian Hyperparameter Optimization for Keras [Video]
   [Notebook]
- Part 8.5: Current Semester's Kaggle [Video] [Notebook]

## Google CoLab Instructions

The following code ensures that Google CoLab is running the correct version of TensorFlow.

Note: using Google CoLab

# Part 8.4: Bayesian Hyperparameter Optimization for Keras

Bayesian Hyperparameter Optimization is a method of finding hyperparameters more efficiently than a grid search. Because each candidate set of hyperparameters requires a retraining of the neural network, it is best to keep the number of candidate sets to a minimum. Bayesian Hyperparameter Optimization achieves this by training a model to predict good candidate sets of hyperparameters. [Cite:snoek2012practical]

- bayesian-optimization
- hyperopt
- spearmint

```
In [2]: # Ignore useless W0819 warnings generated by TensorFlow 2.0.
        # Hopefully can remove this ignore in the future.
        # See https://github.com/tensorflow/tensorflow/issues/31308
        import logging, os
        logging.disable(logging.WARNING)
        os.environ["TF CPP MIN LOG LEVEL"] = "3"
        import pandas as pd
        from scipy.stats import zscore
        # Read the data set
        df = pd.read csv(
            "https://data.heatonresearch.com/data/t81-558/jh-simple-dataset.csv",
            na values=['NA','?'])
        # Generate dummies for job
        df = pd.concat([df,pd.get dummies(df['job'],prefix="job")],axis=1)
        df.drop('job', axis=1, inplace=True)
        # Generate dummies for area
        df = pd.concat([df,pd.get dummies(df['area'],prefix="area")],axis=1)
        df.drop('area', axis=1, inplace=True)
        # Missing values for income
        med = df['income'].median()
        df['income'] = df['income'].fillna(med)
        # Standardize ranges
        df['income'] = zscore(df['income'])
        df['aspect'] = zscore(df['aspect'])
        df['save rate'] = zscore(df['save rate'])
        df['age'] = zscore(df['age'])
        df['subscriptions'] = zscore(df['subscriptions'])
        # Convert to numpy - Classification
        x columns = df.columns.drop('product').drop('id')
```

```
x = df[x_columns].values
dummies = pd.get_dummies(df['product']) # Classification
products = dummies.columns
y = dummies.values
```

Now that we've preprocessed the data, we can begin the hyperparameter optimization. We start by creating a function that generates the model based on just three parameters. Bayesian optimization works on a vector of numbers, not on a problematic notion like how many layers and neurons are on each layer. To represent this complex neuron structure as a vector, we use several numbers to describe this structure.

- **dropout** The dropout percent for each layer.
- neuronPct What percent of our fixed 5,000 maximum number of neurons do we wish to use? This parameter specifies the total count of neurons in the entire network.
- neuronShrink Neural networks usually start with more neurons on the first hidden layer and then decrease this count for additional layers. This percent specifies how much to shrink subsequent layers based on the previous layer.
   We stop adding more layers once we run out of neurons (the count specified by neuronPct).

These three numbers define the structure of the neural network. The commends in the below code show exactly how the program constructs the network.

```
In [3]: import pandas as pd
        import os
        import numpy as np
        import time
        import tensorflow.keras.
        import statistics
        import tensorflow.keras
        from sklearn import metrics
        from sklearn.model selection import StratifiedKFold
        from tensorflow.keras.models import Sequential
        from tensorflow.keras.layers import Dense, Activation, Dropout, InputLayer
        from tensorflow.keras import regularizers
        from tensorflow.keras.callbacks import EarlyStopping
        from sklearn.model selection import StratifiedShuffleSplit
        from sklearn.model selection import ShuffleSplit
        from tensorflow.keras.layers import LeakyReLU,PReLU
        from tensorflow.keras.optimizers import Adam
        def generate model(dropout, neuronPct, neuronShrink):
            # We start with some percent of 5000 starting neurons on
            # the first hidden layer.
            neuronCount = int(neuronPct * 5000)
            # Construct neural network
            model = Sequential()
```

```
# So long as there would have been at least 25 neurons and
# fewer than 10
# layers, create a new layer.
layer = 0
while neuronCount>25 and layer<10:</pre>
    # The first (0th) layer needs an input input dim(neuronCount)
    if layer==0:
        model.add(Dense(neuronCount,
            input dim=x.shape[1],
            activation=PReLU()))
    else:
        model.add(Dense(neuronCount, activation=PReLU()))
    layer += 1
    # Add dropout after each hidden layer
    model.add(Dropout(dropout))
    # Shrink neuron count for each layer
    neuronCount = neuronCount * neuronShrink
model.add(Dense(y.shape[1],activation='softmax')) # Output
return model
```

We can test this code to see how it creates a neural network based on three such parameters.

```
In [4]: # Generate a model and see what the resulting structure looks like.
model = generate_model(dropout=0.2, neuronPct=0.1, neuronShrink=0.25)
model.summary()
```

Model: "sequential"

Layer (type)	Output Shape	Param #
dense (Dense)	(None, 500)	24500
dropout (Dropout)	(None, 500)	Θ
dense_1 (Dense)	(None, 125)	62750
dropout_1 (Dropout)	(None, 125)	0
dense_2 (Dense)	(None, 31)	3937
dropout_2 (Dropout)	(None, 31)	0
dense_3 (Dense)	(None, 7)	224

\_\_\_\_\_\_

Total params: 91,411 Trainable params: 91,411 Non-trainable params: 0 We will now create a function to evaluate the neural network using three such parameters. We use bootstrapping because one training run might have "bad luck" with the assigned random weights. We use this function to train and then evaluate the neural network.

```
In [5]: SPLITS = 2
        EPOCHS = 500
        PATTENCE = 10
        def evaluate network(dropout,learning rate,neuronPct,neuronShrink):
            # Bootstrap
            # for Classification
            boot = StratifiedShuffleSplit(n splits=SPLITS, test size=0.1)
            # for Regression
            # boot = ShuffleSplit(n splits=SPLITS, test size=0.1)
            # Track progress
            mean benchmark = []
            epochs needed = []
            num = 0
            # Loop through samples
            for train, test in boot.split(x,df['product']):
                 start time = time.time()
                 num+=1
                # Split train and test
                x train = x[train]
                y train = y[train]
                x \text{ test} = x[\text{test}]
                y \text{ test} = y[\text{test}]
                 model = generate model(dropout, neuronPct, neuronShrink)
                 model.compile(loss='categorical crossentropy',
                               optimizer=Adam(learning rate=learning rate))
                 monitor = EarlyStopping(monitor='val loss', min delta=1e-3,
                 patience=PATIENCE, verbose=0, mode='auto',
                                          restore best weights=True)
                 # Train on the bootstrap sample
                model.fit(x train,y train,validation data=(x test,y test),
                           callbacks=[monitor], verbose=0, epochs=EPOCHS)
                 epochs = monitor.stopped epoch
                 epochs needed.append(epochs)
                 # Predict on the out of boot (validation)
                 pred = model.predict(x test)
                 # Measure this bootstrap's log loss
                y_compare = np.argmax(y_test,axis=1) # For log loss calculation
                 score = metrics.log loss(y compare, pred)
                 mean benchmark.append(score)
                m1 = statistics.mean(mean benchmark)
```

```
m2 = statistics.mean(epochs_needed)
mdev = statistics.pstdev(mean_benchmark)

# Record this iteration
time_took = time.time() - start_time

tensorflow.keras.backend.clear_session()
return (-m1)
```

You can try any combination of our three hyperparameters, plus the learning rate, to see how effective these four numbers are. Of course, our goal is not to manually choose different combinations of these four hyperparameters; we seek to automate.

-0.6668764846259546

First, we must install the Bayesian optimization package if we are in Colab.

```
In [7]: # HIDE OUTPUT
!pip install bayesian-optimization
```

Requirement already satisfied: bayesian-optimization in /usr/local/lib/pytho n3.7/dist-packages (1.2.0)

Requirement already satisfied: scipy>=0.14.0 in /usr/local/lib/python3.7/dist-packages (from bayesian-optimization) (1.4.1)

Requirement already satisfied: scikit-learn>=0.18.0 in /usr/local/lib/python 3.7/dist-packages (from bayesian-optimization) (1.0.2)

Requirement already satisfied: numpy>=1.9.0 in /usr/local/lib/python3.7/dist -packages (from bayesian-optimization) (1.21.5)

Requirement already satisfied: threadpoolctl>=2.0.0 in /usr/local/lib/python 3.7/dist-packages (from scikit-learn>=0.18.0->bayesian-optimization) (3.1.0) Requirement already satisfied: joblib>=0.11 in /usr/local/lib/python3.7/dist-packages (from scikit-learn>=0.18.0->bayesian-optimization) (1.1.0)

We will now automate this process. We define the bounds for each of these four hyperparameters and begin the Bayesian optimization. Once the program finishes, the best combination of hyperparameters found is displayed. The **optimize** function accepts two parameters that will significantly impact how long the process takes to complete:

- **n\_iter** How many steps of Bayesian optimization that you want to perform. The more steps, the more likely you will find a reasonable maximum.
- **init\_points**: How many steps of random exploration that you want to perform. Random exploration can help by diversifying the exploration space.

```
In [8]: from bayes opt import BayesianOptimization
        import time
        # Supress NaN warnings
        import warnings
        warnings.filterwarnings("ignore", category =RuntimeWarning)
        # Bounded region of parameter space
        pbounds = {'dropout': (0.0, 0.499),
                   'learning rate': (0.0, 0.1),
                   'neuronPct': (0.01, 1),
                   'neuronShrink': (0.01, 1)
        optimizer = BayesianOptimization(
            f=evaluate network,
            pbounds=pbounds,
            verbose=2, # verbose = 1 prints only when a maximum
            # is observed, verbose = 0 is silent
            random state=1,
        start time = time.time()
        optimizer.maximize(init_points=10, n_iter=20,)
        time_took = time.time() - start_time
        print(f"Total runtime: {hms string(time took)}")
        print(optimizer.max)
```

iter	target	dropout	learni	neuronPct	neuron
1	-0.8092	0.2081	0.07203	0.01011	0.3093
2	-0.7167	0.07323	0.009234	0.1944	0.3521
3	-17.87	0.198	0.05388	0.425	0.6884
4	-0.8022	0.102	0.08781	0.03711	0.6738
5	-0.9209	0.2082	0.05587	0.149	0.2061
6	-17.96	0.3996	0.09683	0.3203	0.6954
7	-4.223	0.4373	0.08946	0.09419	0.04866
8	-0.7025	0.08475	0.08781	0.1074	0.4269
9	-8.666	0.478	0.05332	0.695	0.3224
10	-9.785	0.3426	0.08346	0.02811	0.7526
11	-4.881	0.0	0.0	0.01	0.01
12	-21.59	0.2208	0.04135	0.5523	0.7468
13	-1.819	0.0	0.0	1.0	0.01
14	-33.33	0.01058	0.08079	0.9652	0.7051
15	-1.418	0.4963	0.02476	0.9744	0.01896
16	-1.876	0.1247	0.0	0.5781	0.01
17	-1.898	0.0	0.0	0.01	1.0
18	-17.7	0.1621	0.06358	0.4065	0.754
19	-2.674	0.0	0.0	0.01	0.5464
20	-1.931	0.499	0.0	0.5538	0.01
21	-3.402	0.004722	0.05502	0.1704	0.483
22	-2.8	0.08639	0.0838	0.04668	0.6864
23	-20.98	0.1168	0.0447	0.5546	0.9497
24	-4.565	0.2554	0.1	0.8418	0.01
25	-4.724	0.0	0.1	0.3368	0.01
26	-0.6956	0.2505	0.007623	0.01265	0.523
27	-0.7139	0.2967	0.01162	0.3735	0.01708
28	-2.145	0.499	0.0	0.3053	0.2207
29	-2.069	0.0	0.0	0.4808	0.2473
j 30 	-0.7155	0.4082	0.01635	0.02488	0.1694

Total runtime: 1:36:11.56

 ${\text{'target': -0.6955536706512794, 'params': {'dropout': 0.2504561773412203, 'learning_rate': 0.0076232346709142924, 'neuronPct': 0.012648791521811826, 'neuronShrink': 0.5229748831552032}}$ 

As you can see, the algorithm performed 30 total iterations. This total iteration count includes ten random and 20 optimization iterations.



# T81-558: Applications of Deep Neural Networks

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- Part 8.5: Current Semester's Kaggle [Video] [Notebook]

# Google CoLab Instructions

The following code ensures that Google CoLab is running the correct version of TensorFlow.

Note: using Google CoLab

# Part 8.5: Current Semester's Kaggle

Kaggke competition site for current semester:

Spring 2023 Kaggle Assignment

Previous Kaggle competition sites for this class (NOT this semester's assignment, feel free to use code):

- Fall 2022 Kaggle Assignment
- Spring 2022 Kaggle Assignment
- Fall 2021 Kaggle Assignment
- Spring 2021 Kaggle Assignment
- Fall 2020 Kaggle Assignment
- Spring 2020 Kaggle Assignment
- Fall 2019 Kaggle Assignment
- Spring 2019 Kaggle Assignment
- Fall 2018 Kaggle Assignment
- Spring 2018 Kaggle Assignment
- Fall 2017 Kaggle Assignment
- Spring 2017 Kaggle Assignment
- Fall 2016 Kaggle Assignment

#### Iris as a Kaggle Competition

If I used the Iris data as a Kaggle, I would give you the following three files:

- kaggle iris test.csv The data that Kaggle will evaluate you on. It contains only input; you must provide answers. (contains x)
- kaggle iris train.csv The data that you will use to train. (contains x and y)
- kaggle iris sample.csv A sample submission for Kaggle. (contains x and y)

Important features of the Kaggle iris files (that differ from how we've previously seen files):

- The iris species is already index encoded.
- Your training data is in a separate file.
- You will load the test data to generate a submission file.

The following program generates a submission file for "Iris Kaggle". You can use it as a starting point for assignment 3.

```
from sklearn.model selection import train test split
import tensorflow as tf
import numpy as np
from tensorflow.keras.models import Sequential
from tensorflow.keras.layers import Dense, Activation
from tensorflow.keras.callbacks import EarlyStopping
df train = pd.read csv(
    "https://data.heatonresearch.com/data/t81-558/datasets/"+\
    "kaggle iris train.csv", na values=['NA','?'])
# Encode feature vector
df train.drop('id', axis=1, inplace=True)
num classes = len(df train.groupby('species').species.nunique())
print("Number of classes: {}".format(num classes))
# Convert to numpy - Classification
x = df_train[['sepal_l', 'sepal_w', 'petal_l', 'petal_w']].values
dummies = pd.get dummies(df train['species']) # Classification
species = dummies.columns
y = dummies.values
# Split into train/test
x train, x test, y train, y test = train test split(
    x, y, test size=0.25, random state=45)
# Train, with early stopping
model = Sequential()
model.add(Dense(50, input dim=x.shape[1], activation='relu'))
model.add(Dense(25))
model.add(Dense(y.shape[1],activation='softmax'))
model.compile(loss='categorical crossentropy', optimizer='adam')
monitor = EarlyStopping(monitor='val loss', min delta=1e-3,
                        patience=5, verbose=1, mode='auto',
                       restore best weights=True)
model.fit(x train,y train,validation data=(x test,y test),
          callbacks=[monitor], verbose=0, epochs=1000)
```

```
Number of classes: 3
Restoring model weights from the end of the best epoch: 103.
Epoch 108: early stopping
```

Out[2]: <keras.callbacks.History at 0x7f05e7452710>

Now that we've trained the neural network, we can check its log loss.

```
In [3]: from sklearn import metrics

# Calculate multi log loss error
pred = model.predict(x_test)
score = metrics.log_loss(y_test, pred)
print("Log loss score: {}".format(score))
```

```
Log loss score: 0.10988010508939623
```

Now we are ready to generate the Kaggle submission file. We will use the iris test data that does not contain a y target value. It is our job to predict this value and submit it to Kaggle.

```
In [4]: # Generate Kaggle submit file
        # Encode feature vector
        df test = pd.read csv(
            "https://data.heatonresearch.com/data/t81-558/datasets/"+\
            "kaggle iris test.csv", na values=['NA','?'])
        # Convert to numpy - Classification
        ids = df test['id']
        df test.drop('id', axis=1, inplace=True)
        x = df_test[['sepal_l', 'sepal_w', 'petal_l', 'petal_w']].values
        y = dummies.values
        # Generate predictions
        pred = model.predict(x)
        #pred
        # Create submission data set
        df submit = pd.DataFrame(pred)
        df submit.insert(0,'id',ids)
        df submit.columns = ['id','species-0','species-1','species-2']
        # Write submit file locally
        df submit.to csv("iris submit.csv", index=False)
        print(df_submit[:5])
          id species-0 species-1 species-2
       0 100 0.022300 0.777859 0.199841
```

```
id species-0 species-1 species-2
0 100 0.022300 0.777859 0.199841
1 101 0.001309 0.273849 0.724842
2 102 0.001153 0.319349 0.679498
3 103 0.958006 0.041989 0.000005
4 104 0.976932 0.023066 0.000002
```

## MPG as a Kaggle Competition (Regression)

If the Auto MPG data were used as a Kaggle, you would be given the following three files:

- kaggle\_mpg\_test.csv The data that Kaggle will evaluate you on. Contains only input, you must provide answers. (contains x)
- kaggle mpg train.csv The data that you will use to train. (contains x and y)
- kaggle mpg sample.csv A sample submission for Kaggle. (contains x and y)

Important features of the Kaggle iris files (that differ from how we've previously seen files):

The following program generates a submission file for "MPG Kaggle".

```
In [5]: # HIDE OUTPUT
        from tensorflow.keras.models import Sequential
        from tensorflow.keras.layers import Dense, Activation
        from sklearn.model selection import train test split
        from tensorflow.keras.callbacks import EarlyStopping
        import pandas as pd
        import io
        import os
        import requests
        import numpy as np
        from sklearn import metrics
        save path = "."
        df = pd.read csv(
            "https://data.heatonresearch.com/data/t81-558/datasets/"+\
            "kaggle auto train.csv",
            na values=['NA', '?'])
        cars = df['name']
        # Handle missing value
        df['horsepower'] = df['horsepower'].fillna(df['horsepower'].median())
        # Pandas to Numpy
        x = df[['cylinders', 'displacement', 'horsepower', 'weight',
               'acceleration', 'year', 'origin']].values
        y = df['mpg'].values # regression
        # Split into train/test
        x train, x test, y train, y test = train test split(
            x, y, test size=0.25, random state=42)
        # Build the neural network
        model = Sequential()
        model.add(Dense(25, input dim=x.shape[1], activation='relu')) # Hidden 1
        model.add(Dense(10, activation='relu')) # Hidden 2
        model.add(Dense(1)) # Output
        model.compile(loss='mean squared error', optimizer='adam')
        monitor = EarlyStopping(monitor='val loss', min delta=1e-3, patience=5,
                                verbose=1, mode='auto', restore best weights=True)
        model.fit(x train,y train,validation data=(x test,y test),
                  verbose=2, callbacks=[monitor], epochs=1000)
        # Predict
        pred = model.predict(x test)
```

```
Epoch 1/1000
9/9 - 1s - loss: 1797.5945 - val loss: 1272.4421 - 1s/epoch - 144ms/step
Epoch 2/1000
9/9 - 0s - loss: 574.7726 - val loss: 734.3082 - 92ms/epoch - 10ms/step
Epoch 3/1000
9/9 - 0s - loss: 487.3118 - val loss: 446.3558 - 76ms/epoch - 8ms/step
Epoch 4/1000
9/9 - 0s - loss: 326.7128 - val loss: 321.7191 - 96ms/epoch - 11ms/step
Epoch 5/1000
9/9 - 0s - loss: 294.8217 - val_loss: 271.3473 - 70ms/epoch - 8ms/step
Epoch 6/1000
9/9 - 0s - loss: 259.8376 - val loss: 239.6796 - 116ms/epoch - 13ms/step
Epoch 7/1000
9/9 - 0s - loss: 250.4708 - val loss: 227.4295 - 73ms/epoch - 8ms/step
Epoch 8/1000
9/9 - 0s - loss: 227.1252 - val loss: 198.4167 - 125ms/epoch - 14ms/step
Epoch 9/1000
9/9 - 0s - loss: 225.6681 - val loss: 195.5055 - 95ms/epoch - 11ms/step
Epoch 10/1000
9/9 - 0s - loss: 209.1198 - val loss: 184.1092 - 121ms/epoch - 13ms/step
Epoch 11/1000
9/9 - 0s - loss: 195.4801 - val loss: 176.0311 - 108ms/epoch - 12ms/step
Epoch 12/1000
9/9 - 0s - loss: 198.6493 - val loss: 168.1613 - 163ms/epoch - 18ms/step
Epoch 13/1000
9/9 - 0s - loss: 198.5606 - val loss: 196.0306 - 114ms/epoch - 13ms/step
Epoch 14/1000
9/9 - 0s - loss: 184.3067 - val loss: 179.8450 - 99ms/epoch - 11ms/step
Epoch 15/1000
9/9 - 0s - loss: 178.6627 - val loss: 148.1014 - 80ms/epoch - 9ms/step
Epoch 16/1000
9/9 - 0s - loss: 154.0201 - val loss: 129.9253 - 74ms/epoch - 8ms/step
Epoch 17/1000
9/9 - 0s - loss: 145.2373 - val loss: 124.0609 - 79ms/epoch - 9ms/step
Epoch 18/1000
9/9 - 0s - loss: 140.0318 - val loss: 116.7844 - 86ms/epoch - 10ms/step
Epoch 19/1000
9/9 - 0s - loss: 135.1688 - val loss: 115.0745 - 136ms/epoch - 15ms/step
Epoch 20/1000
9/9 - 0s - loss: 132.8391 - val_loss: 106.9831 - 169ms/epoch - 19ms/step
Epoch 21/1000
9/9 - 0s - loss: 123.6673 - val loss: 105.7211 - 95ms/epoch - 11ms/step
Epoch 22/1000
9/9 - 0s - loss: 123.7169 - val_loss: 99.6713 - 112ms/epoch - 12ms/step
Epoch 23/1000
9/9 - 0s - loss: 118.0815 - val loss: 96.0683 - 150ms/epoch - 17ms/step
Epoch 24/1000
9/9 - 0s - loss: 114.6363 - val_loss: 99.1486 - 153ms/epoch - 17ms/step
Epoch 25/1000
9/9 - 0s - loss: 112.3965 - val loss: 93.8642 - 180ms/epoch - 20ms/step
Epoch 26/1000
9/9 - 0s - loss: 111.2470 - val_loss: 88.3417 - 139ms/epoch - 15ms/step
Epoch 27/1000
9/9 - 0s - loss: 107.8639 - val loss: 86.7927 - 135ms/epoch - 15ms/step
Epoch 28/1000
9/9 - 0s - loss: 103.0426 - val loss: 89.0441 - 101ms/epoch - 11ms/step
```

```
Epoch 29/1000
9/9 - 0s - loss: 110.6277 - val loss: 82.4294 - 159ms/epoch - 18ms/step
Epoch 30/1000
9/9 - 0s - loss: 100.3681 - val loss: 90.8037 - 82ms/epoch - 9ms/step
Epoch 31/1000
9/9 - 0s - loss: 105.4711 - val loss: 79.2106 - 76ms/epoch - 8ms/step
Epoch 32/1000
9/9 - 0s - loss: 98.7603 - val loss: 79.9620 - 73ms/epoch - 8ms/step
Epoch 33/1000
9/9 - 0s - loss: 94.7678 - val loss: 76.8616 - 78ms/epoch - 9ms/step
Epoch 34/1000
9/9 - 0s - loss: 93.8199 - val loss: 77.0823 - 76ms/epoch - 8ms/step
Epoch 35/1000
9/9 - 0s - loss: 94.8746 - val loss: 73.9967 - 62ms/epoch - 7ms/step
Epoch 36/1000
9/9 - 0s - loss: 95.3178 - val_loss: 73.0059 - 60ms/epoch - 7ms/step
Epoch 37/1000
9/9 - 0s - loss: 91.1315 - val loss: 80.8389 - 57ms/epoch - 6ms/step
Epoch 38/1000
9/9 - 0s - loss: 96.4810 - val loss: 77.8854 - 59ms/epoch - 7ms/step
Epoch 39/1000
9/9 - 0s - loss: 91.1039 - val loss: 69.9539 - 40ms/epoch - 4ms/step
Epoch 40/1000
9/9 - 0s - loss: 86.9596 - val loss: 69.3511 - 43ms/epoch - 5ms/step
Epoch 41/1000
9/9 - 0s - loss: 87.6142 - val loss: 70.1390 - 57ms/epoch - 6ms/step
Epoch 42/1000
9/9 - 0s - loss: 88.0185 - val loss: 73.6168 - 38ms/epoch - 4ms/step
Epoch 43/1000
9/9 - 0s - loss: 92.8655 - val loss: 67.5213 - 38ms/epoch - 4ms/step
Epoch 44/1000
9/9 - 0s - loss: 88.5278 - val loss: 69.9708 - 59ms/epoch - 7ms/step
Epoch 45/1000
9/9 - 0s - loss: 82.9339 - val loss: 70.3786 - 39ms/epoch - 4ms/step
Epoch 46/1000
9/9 - 0s - loss: 81.7092 - val loss: 63.3550 - 59ms/epoch - 7ms/step
Epoch 47/1000
9/9 - 0s - loss: 81.1514 - val_loss: 78.7681 - 59ms/epoch - 7ms/step
Epoch 48/1000
9/9 - 0s - loss: 99.3562 - val loss: 62.8894 - 64ms/epoch - 7ms/step
Epoch 49/1000
9/9 - 0s - loss: 96.8292 - val loss: 67.8047 - 55ms/epoch - 6ms/step
Epoch 50/1000
9/9 - 0s - loss: 88.7995 - val loss: 67.5249 - 56ms/epoch - 6ms/step
Epoch 51/1000
9/9 - 0s - loss: 80.6064 - val loss: 96.2975 - 58ms/epoch - 6ms/step
Epoch 52/1000
9/9 - 0s - loss: 95.2732 - val loss: 62.4323 - 39ms/epoch - 4ms/step
Epoch 53/1000
9/9 - 0s - loss: 75.1992 - val loss: 64.0174 - 39ms/epoch - 4ms/step
Epoch 54/1000
9/9 - 0s - loss: 75.5173 - val loss: 57.8594 - 40ms/epoch - 4ms/step
Epoch 55/1000
9/9 - 0s - loss: 72.6369 - val loss: 56.2216 - 47ms/epoch - 5ms/step
Epoch 56/1000
9/9 - 0s - loss: 72.8636 - val loss: 55.3956 - 54ms/epoch - 6ms/step
```

```
Epoch 57/1000
9/9 - 0s - loss: 69.0251 - val loss: 70.7940 - 56ms/epoch - 6ms/step
Epoch 58/1000
9/9 - 0s - loss: 75.8152 - val loss: 63.7728 - 37ms/epoch - 4ms/step
Epoch 59/1000
9/9 - 0s - loss: 71.6866 - val loss: 59.5908 - 41ms/epoch - 5ms/step
Epoch 60/1000
9/9 - 0s - loss: 69.3349 - val loss: 52.7848 - 38ms/epoch - 4ms/step
Epoch 61/1000
9/9 - 0s - loss: 67.8410 - val loss: 53.5977 - 54ms/epoch - 6ms/step
Epoch 62/1000
9/9 - 0s - loss: 68.4640 - val loss: 53.6664 - 39ms/epoch - 4ms/step
Epoch 63/1000
9/9 - 0s - loss: 63.7229 - val loss: 52.4224 - 44ms/epoch - 5ms/step
Epoch 64/1000
9/9 - 0s - loss: 69.8485 - val loss: 59.1973 - 53ms/epoch - 6ms/step
Epoch 65/1000
9/9 - 0s - loss: 75.7193 - val loss: 70.1342 - 37ms/epoch - 4ms/step
Epoch 66/1000
9/9 - 0s - loss: 87.7418 - val loss: 55.3687 - 38ms/epoch - 4ms/step
Epoch 67/1000
9/9 - 0s - loss: 72.8599 - val loss: 52.9028 - 44ms/epoch - 5ms/step
Epoch 68/1000
9/9 - 0s - loss: 69.9528 - val loss: 49.9109 - 38ms/epoch - 4ms/step
Epoch 69/1000
9/9 - 0s - loss: 62.7782 - val loss: 46.6361 - 39ms/epoch - 4ms/step
Epoch 70/1000
9/9 - 0s - loss: 58.4024 - val loss: 50.8190 - 38ms/epoch - 4ms/step
Epoch 71/1000
9/9 - 0s - loss: 63.5687 - val loss: 46.6161 - 44ms/epoch - 5ms/step
Epoch 72/1000
9/9 - 0s - loss: 65.9290 - val loss: 47.1278 - 40ms/epoch - 4ms/step
Epoch 73/1000
9/9 - 0s - loss: 74.9235 - val loss: 61.1282 - 42ms/epoch - 5ms/step
Epoch 74/1000
9/9 - 0s - loss: 63.6773 - val loss: 45.0233 - 39ms/epoch - 4ms/step
Epoch 75/1000
9/9 - 0s - loss: 55.8287 - val_loss: 59.8986 - 41ms/epoch - 5ms/step
Epoch 76/1000
9/9 - 0s - loss: 58.9969 - val loss: 52.0535 - 39ms/epoch - 4ms/step
Epoch 77/1000
9/9 - 0s - loss: 60.7104 - val loss: 43.0530 - 46ms/epoch - 5ms/step
Epoch 78/1000
9/9 - 0s - loss: 59.7358 - val loss: 45.3669 - 41ms/epoch - 5ms/step
Epoch 79/1000
9/9 - 0s - loss: 60.9792 - val loss: 40.7967 - 41ms/epoch - 5ms/step
Epoch 80/1000
9/9 - 0s - loss: 58.0294 - val loss: 49.0612 - 42ms/epoch - 5ms/step
Epoch 81/1000
9/9 - 0s - loss: 57.6733 - val loss: 41.7604 - 44ms/epoch - 5ms/step
Epoch 82/1000
9/9 - 0s - loss: 50.3309 - val loss: 39.1461 - 38ms/epoch - 4ms/step
Epoch 83/1000
9/9 - 0s - loss: 54.2316 - val loss: 40.8561 - 36ms/epoch - 4ms/step
Epoch 84/1000
9/9 - 0s - loss: 66.4084 - val loss: 38.1869 - 60ms/epoch - 7ms/step
```

```
Epoch 85/1000
9/9 - 0s - loss: 50.0778 - val loss: 37.8852 - 56ms/epoch - 6ms/step
Epoch 86/1000
9/9 - 0s - loss: 47.0763 - val loss: 37.3743 - 39ms/epoch - 4ms/step
Epoch 87/1000
9/9 - 0s - loss: 46.1752 - val loss: 45.8444 - 45ms/epoch - 5ms/step
Epoch 88/1000
9/9 - 0s - loss: 49.4047 - val loss: 37.3778 - 40ms/epoch - 4ms/step
Epoch 89/1000
9/9 - 0s - loss: 46.5478 - val loss: 36.2859 - 38ms/epoch - 4ms/step
Epoch 90/1000
9/9 - 0s - loss: 44.7429 - val loss: 47.2213 - 38ms/epoch - 4ms/step
Epoch 91/1000
9/9 - 0s - loss: 49.7726 - val_loss: 52.5501 - 42ms/epoch - 5ms/step
Epoch 92/1000
9/9 - 0s - loss: 53.5449 - val loss: 62.3078 - 57ms/epoch - 6ms/step
Epoch 93/1000
9/9 - 0s - loss: 54.7558 - val loss: 51.2010 - 43ms/epoch - 5ms/step
Epoch 94/1000
Restoring model weights from the end of the best epoch: 89.
9/9 - 0s - loss: 52.3631 - val_loss: 42.2640 - 69ms/epoch - 8ms/step
Epoch 94: early stopping
```

Now that we've trained the neural network, we can check its RMSE error.

```
In [6]: import numpy as np

# Measure RMSE error. RMSE is common for regression.
score = np.sqrt(metrics.mean_squared_error(pred,y_test))
print("Final score (RMSE): {}".format(score))
```

Final score (RMSE): 6.023776405947501

Now we are ready to generate the Kaggle submission file. We will use the MPG test data that does not contain a y target value. It is our job to predict this value and submit it to Kaggle.

```
In [7]: import pandas as pd

# Generate Kaggle submit file

# Encode feature vector

df_test = pd.read_csv(
        "https://data.heatonresearch.com/data/t81-558/datasets/"+\
        "kaggle_auto_test.csv", na_values=['NA','?'])

# Convert to numpy - regression
ids = df_test['id']
df_test.drop('id', axis=1, inplace=True)

# Handle missing value
df_test['horsepower'] = df_test['horsepower'].\
        fillna(df['horsepower'].median())

x = df_test[['cylinders', 'displacement', 'horsepower', 'weight',
```

```
'acceleration', 'year', 'origin']].values

# Generate predictions
pred = model.predict(x)
#pred

# Create submission data set

df_submit = pd.DataFrame(pred)
df_submit.insert(0,'id',ids)
df_submit.columns = ['id','mpg']

# Write submit file locally
df_submit.to_csv("auto_submit.csv", index=False)
print(df_submit[:5])
```

```
id mpg
0 350 27.158819
1 351 24.450621
2 352 24.913355
3 353 26.994867
4 354 26.669268
```