mlp

May 7, 2022

1 Lab 3: Multi-Layers Perceptron

2 Introduction

The objective of this lab is to dive into particular kind of neural network: the *Multi-Layer Perceptron* (MLP).

To start, let us take the dataset from the previous lab (hydrodynamics of sailing boats) and use scikit-learn to train a MLP instead of our hand-made single perceptron. The code below is already complete and is meant to give you an idea of how to construct an MLP with scikit-learn. You can execute it, taking the time to understand the idea behind each cell.

```
[1]: # Importing the dataset
import numpy as np
dataset = np.genfromtxt("yacht_hydrodynamics.data", delimiter='')
X = dataset[:, :-1]
Y = dataset[:, -1]
```

```
[2]: # Preprocessing: scale input data
from sklearn.preprocessing import StandardScaler
sc = StandardScaler()
X = sc.fit_transform(X)
```

```
[3]: # Split dataset into training and test set

from sklearn.model_selection import train_test_split
x_train, x_test, y_train, y_test = train_test_split(X, Y,random_state=1,___

otest_size = 0.20)
```

```
[4]: # Define a multi-layer perceptron (MLP) network for regression

from sklearn.neural_network import MLPRegressor

mlp = MLPRegressor(max_iter=3000, random_state=1) # define the model, with_
default params

mlp.fit(x_train, y_train) # train the MLP
```

[4]: MLPRegressor(max_iter=3000, random_state=1)

```
[5]: # Evaluate the model from matplotlib import pyplot as plt
```

```
print('Train score: ', mlp.score(x_train, y_train))
print('Test score: ', mlp.score(x_test, y_test))
#plt.plot(mlp.loss_curve_)
#plt.xlabel("Iterations")
#plt.ylabel("Loss")
```

Train score: 0.9940765369322633 Test score: 0.9899773031580283

```
[7]: # Plot the results
   num_samples_to_plot = 20
   #plt.plot(y_test[0:num_samples_to_plot], 'ro', label='y')
   yw = mlp.predict(x_test)
   #plt.plot(yw[0:num_samples_to_plot], 'bx', label='$\hat{y}$')
   #plt.legend()
   #plt.xlabel("Examples")
   #plt.ylabel("f(examples)")
```

2.0.1 Analyzing the network

Many details of the network are currently hidden as default parameters.

Using the documentation of the MLPRegressor, answer the following questions.

- What is the structure of the network? The network is a multi-layer perceptron with hidden layers which we can choose se number and the activation function.
- What it is the algorithm used for training? Is there algorithm available that we mentioned during the courses? The training algorithm is the stochastic gradient descent. Yes it was mentionned during the courses.
- How does the training algorithm decides to stop the training? We have tol that is the minimum desired improvement of the loss or score for consecutive iterations. The training stops when when there is no change.

3 Onto a more challenging dataset: house prices

For the rest of this lab, we will use the (more challenging) California Housing Prices dataset.

```
[7]: # clean all previously defined variables for the sailing boats %reset -f
```

```
[8]: """Import the required modules"""
from sklearn.datasets import fetch_california_housing
import pandas as pd

num_samples = 2000 # only use the first N samples to limit training time

cal_housing = fetch_california_housing()
```

```
X = pd.DataFrame(cal housing.data,columns=cal housing.feature_names)[:
 →num_samples]
Y = cal_housing.target[:num_samples]
X.head(10) # print the first 10 values
```

```
[8]:
       {	t MedInc}
               HouseAge AveRooms
                                   AveBedrms
                                              Population AveOccup Latitude
    0 8.3252
                    41.0
                         6.984127
                                     1.023810
                                                    322.0
                                                          2.555556
                                                                        37.88
    1 8.3014
                    21.0
                         6.238137
                                     0.971880
                                                   2401.0 2.109842
                                                                        37.86
    2 7.2574
                   52.0
                         8.288136
                                    1.073446
                                                    496.0 2.802260
                                                                        37.85
    3 5.6431
                   52.0 5.817352
                                    1.073059
                                                    558.0 2.547945
                                                                        37.85
    4 3.8462
                   52.0 6.281853
                                    1.081081
                                                   565.0 2.181467
                                                                        37.85
    5 4.0368
                   52.0 4.761658
                                     1.103627
                                                   413.0 2.139896
                                                                       37.85
                   52.0 4.931907
    6 3.6591
                                    0.951362
                                                   1094.0 2.128405
                                                                        37.84
    7 3.1200
                   52.0 4.797527
                                                                        37.84
                                     1.061824
                                                   1157.0 1.788253
    8 2.0804
                   42.0 4.294118
                                                   1206.0 2.026891
                                                                        37.84
                                     1.117647
    9 3.6912
                   52.0 4.970588
                                     0.990196
                                                   1551.0 2.172269
                                                                        37.84
```

Longitude

- -122.23 0
- -122.221
- 2 -122.24
- 3 -122.25
- -122.254
- 5 -122.25
- 6 -122.25
- 7
- -122.258 -122.26
- 9 -122.25

Note that each row of the dataset represents a **group of houses** (one district). The target variable denotes the average house value in units of 100.000 USD. Median Income is per 10.000 USD.

3.0.1 Extracting a subpart of the dataset for testing

• Split the dataset between a training set (75%) and a test set (25%)

Please use the conventional names X_train, X_test, y_train and y_test.

```
[9]: # TODO
     from sklearn.model_selection import train_test_split
     x_train, x_test, y_train, y_test = train_test_split(X, Y,random_state=1,__
      \rightarrowtest size = 0.25)
```

3.0.2 Scaling the input data

A step of scaling of the data is often useful to ensure that all input data centered on 0 and with a fixed variance.

Standardization of a dataset is a common requirement for many machine learning estimators: they might behave badly if the individual features do not more or less look like standard normally distributed data (e.g. Gaussian with 0 mean and unit variance). The function StandardScaler from sklearn.preprocessing computes the standard score of a sample as:

```
z = (x - u) / s
```

where u is the mean of the training samples, and s is the standard deviation of the training samples.

Centering and scaling happen independently on each feature by computing the relevant statistics on the samples in the training set. Mean and standard deviation are then stored to be used on later data using transform.

- Apply the standard scaler to both the training dataset (X_train) and the test dataset (X_test).
- Make sure that **exactly the same transformation** is applied to both datasets.

Documentation of standard scaler in scikit learn

```
[10]: # TODO
from sklearn.preprocessing import StandardScaler

sc = StandardScaler()
X_train = sc.fit_transform(x_train)
X_test = sc.fit_transform(x_test)
```

3.1 Overfitting

In this part, we are only interested in maximizing the **train score**, i.e., having the network memorize the training examples as well as possible.

• Propose a parameterization of the network (shape and learning parameters) that will maximize the train score (without considering the test score).

While doing this, you should (1) remain within two minutes of training time, and (2) obtain a score that is greater than 0.90.

• Is the **test** score substantially smaller than the **train** score (indicator of overfitting)?

Yes because with a good train score we tend to overfit

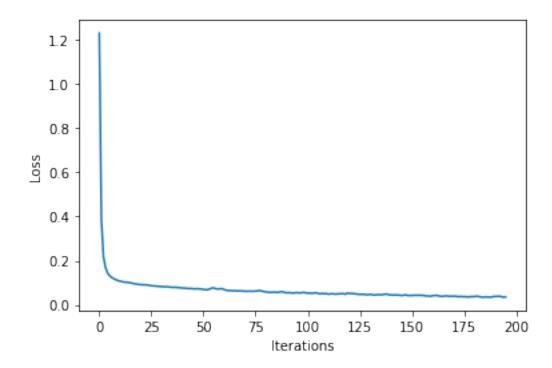
• Explain how the parameters you chose allow the learned model to overfit.

To increase the train score, we can start by adding mode hidden layers to predicate better the training value of the houses and delete the validation fraction so that all of the set is used for training. We can also decrease the tolerance, increase the number of iterations of the algorithm to get a more precise results.

```
[11]: # TODO
from sklearn.neural_network import MLPRegressor
from matplotlib import pyplot as plt
```

Train score: 0.9143311098389366 Test score: 0.3194622548310331

[11]: Text(0, 0.5, 'Loss')



3.2 Hyperparameter tuning

In this section, we are now interested in maximizing the ability of the network to predict the value of unseen examples, i.e., maximizing the **test** score. You should experiment with the possible parameters of the network in order to obtain a good test score, ideally with a small learning time.

Parameters to vary:

- number and size of the hidden layers
- activation function
- stopping conditions
- maximum number of iterations
- initial learning rate value

Results to present for the tested configurations:

- Train/test score
- training time

Present in a table the various parameters tested and the associated results. You can find in the last cell of the notebook a code snippet that will allow you to plot tables from python structure. Be methodical in the way your run your experiments and collect data. For each run, you should record the parameters and results into an external data structure.

(Note that, while we encourage you to explore the solution space manually, there are existing methods in scikit-learn and other learning framework to automate this step as well, e.g., GridSearchCV)

```
/usr/local/insa/anaconda/lib/python3.8/site-
packages/sklearn/neural_network/_multilayer_perceptron.py:614:
ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't converged yet.
warnings.warn(
```

[12]: MLPRegressor()

```
[13]: conf1 = {"hidden_layer_sizes" : (100,100,100, 100), "activation" : 'relu', [13]
       →"tol" :0.0001, "early_stopping" : False, "max_iter":200, □

¬"learning_rate_init":0.001}
      test_conf(conf1)
[13]: MLPRegressor(hidden_layer_sizes=(100, 100, 100, 100))
[14]: conf2 = {"hidden_layer_sizes" : (10,), "activation" : 'relu', "tol" :0.0001, __
      → "early_stopping": False, "max_iter":200, "learning_rate_init":0.001}
      test conf(conf2)
     /usr/local/insa/anaconda/lib/python3.8/site-
     packages/sklearn/neural_network/_multilayer_perceptron.py:614:
     ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and
     the optimization hasn't converged yet.
       warnings.warn(
[14]: MLPRegressor(hidden_layer_sizes=(10,))
[15]: conf3 = {"hidden_layer_sizes" : (100,), "activation" : 'identity', "tol" :0.
      →0001, "early_stopping" : False, "max_iter":200, "learning_rate_init":0.001}
      test conf(conf3)
[15]: MLPRegressor(activation='identity')
[16]: conf4 = {"hidden_layer_sizes" : (100,), "activation" : 'logistic', "tol" :0.
      →0001, "early_stopping" : False, "max_iter":200, "learning_rate_init":0.001}
      test conf(conf4)
     /usr/local/insa/anaconda/lib/python3.8/site-
     packages/sklearn/neural_network/_multilayer_perceptron.py:614:
     ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and
     the optimization hasn't converged yet.
       warnings.warn(
[16]: MLPRegressor(activation='logistic')
[17]: conf5 = {"hidden_layer_sizes": (100,), "activation": 'tanh', "tol": 0.0001, |
       →"early_stopping" : False, "max_iter":200, "learning_rate_init":0.001}
      test conf(conf5)
     /usr/local/insa/anaconda/lib/python3.8/site-
     packages/sklearn/neural_network/_multilayer_perceptron.py:614:
     ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and
     the optimization hasn't converged yet.
       warnings.warn(
```

```
[17]: MLPRegressor(activation='tanh')
[18]: conf6 = {"hidden_layer_sizes" : (100,), "activation" : 'relu', "tol" :0.01,
       →"early_stopping" : False, "max_iter":200, "learning_rate_init":0.001}
      test conf(conf6)
[18]: MLPRegressor(tol=0.01)
[19]: conf7 = {"hidden_layer_sizes": (100,), "activation": 'relu', "tol": 0.0001, |
       → "early_stopping": True, "max_iter":200, "learning_rate_init":0.001}
      test conf(conf7)
[19]: MLPRegressor(early_stopping=True)
[20]: conf8 = {"hidden layer sizes": (100,), "activation": 'relu', "tol": 0.0001, |
       → "early_stopping": False, "max_iter":2000, "learning_rate_init":0.001}
      test conf(conf8)
[20]: MLPRegressor(max_iter=2000)
[21]: conf9 = {"hidden_layer_sizes" : (100,), "activation" : 'relu', "tol" :0.0001,
      -- "early_stopping" : False, "max_iter":200, "learning_rate_init":0.1}
      final_mlp = test_conf(conf9)
[22]: #The best configuration is one that gives us the best score for both training.
       \rightarrow and testing set
      #We should then compromise on the number of layers and the maximum of iterations
      best_conf = {"hidden_layer_sizes" : (100,25), "activation" : 'relu', "tol" :0.
       →0001, "early_stopping" : True,
                   "learning_rate" : 'adaptive', "validation_fraction" : 0.1, __
       →"max_iter":500, "learning_rate_init":0.001 }
      test_conf(best_conf)
[22]: MLPRegressor(early_stopping=True, hidden_layer_sizes=(100, 25),
                   learning_rate='adaptive', max_iter=500)
[23]: import numpy as np
      table = pd.DataFrame.from_dict(data)
      table = table.replace(np.nan, '-')
      table = table.sort_values(by='Test score', ascending=False)
      table
[23]:
            hidden_layer_sizes activation
                                                   early_stopping max_iter \
                                              tol
      10
                     (100, 25)
                                     relu 0.0001
                                                             True
                                                                        500
      7
```

True

200

relu 0.0001

(100,)

0		(100,) relu	0.0001	False	200	
6		(100,) relu	0.0100	False	200	
9		(100,) relu	0.0001	False	200	
5		(100,) tanh	0.0001	False	200	
4		(100,) logistic	0.0001	False	200	
8		(100,) relu	0.0001	False	2000	
3	(100,) identity	0.0001	False	200	
2	(10,) relu	0.0001	False	200	
1	(100, 100,	100, 100) relu	0.0001	False	200	
	learning_r	ate init	Time	Train score	Test score l	Learning_rate	\
10	1001111116_1	0.001	13.593365	0.800659	0.796612	adaptive	`
7		0.001	18.896147	0.797888	0.785586	- daap0110	
0		0.001	25.887275	0.818538	0.781142	_	
6		0.001	2.804348	0.739251	0.767593	_	
9		0.100	4.988372	0.822721	0.765393	_	
5		0.001	35.992563	0.788781	0.753041	_	
4		0.001	28.594653	0.740026	0.752026	_	
8		0.001	48.092507	0.838026	0.737605	_	
3		0.001	4.758420	0.688997	0.725396	-	
2		0.001	0.832452	0.736469	0.687174	_	
1		0.001	36.901365	0.902037	0.566535	_	
validation_fraction							
10	varraaoron_	0.1					
7		-					
0		_					
6		_					
9		_					
5		_					
4		_					
8		_					
3		_					

3.3 Evaluation

2

- From your experiments, what seems to be the best model (i.e. set of parameters) for predicting the value of a house?
 - Influences of the parameters :
 - * the higher the number and size of hidden layer the smaller the test score, the best the train score and the higher the computation time
 - * the relu activation function give the best test and train score, followed by the logistic activation function but that one takes more time to compute
 - * less iterations give the best test score, worst train score and best computation time
 - * early_stopping set to True is good to optimize the computation time

- * higher tolerance gives the quickest computation time but decreases the test and train score
- * higher initial learning rate increases the train score but decreases the test score
- * validation fraction helps with overfitting so it improves the test score
- * learning rate optimize time
- Our best config would be : {"hidden_layer_sizes" : (100,25), "activation" : 'relu', "tol"
 :0.0001, "early_stopping" : True, "learning_rate" : 'adaptive', "validation_fraction" : 0.1, "max_iter":500, "learning_rate_init":0.001 }

Unless you used cross-validation, you have probably used the "test" set to select the best model among the ones you experimented with. Since your model is the one that worked best on the "test" set, your selection is *biased*.

In all rigor the original dataset should be split in three:

- the **training set**, on which each model is trained
- the validation set, that is used to pick the best parameters of the model
- the **test set**, on which we evaluate the final model

Evaluate the score of your algorithm on a test set that was not used for training nor for model selection.

```
Test score 1= -9.00692735001358
Test score 2= -0.3160968532235562
Test score 3= -0.10173092242689696
```

We can see that with data outside of our training or testing set, the test score are terrible. This is because our algorithm is biased and our training set is small.