# Digit recognizer: Model comparison and hyperparameter tuning

This project aims to develop a machine learning model for recognising handwritten digits and is part of the Kaggle competition Digit Recognizer. The MNIST dataset, the de facto "hello world" dataset of computer vision, is used to train and test the models.

We compare different machine learning models: logistic regression, fully connected neural network, and convolutional neural network. We also perform hyperparameter tuning to achieve the best performance.

```
In [1]:
         import matplotlib as mpl
         import matplotlib.pyplot as plt
         import math
         import numpy as np
         import pandas as pd
         import seaborn as sns
         from sklearn.linear model import LogisticRegression
         from sklearn.metrics import accuracy score, confusion matrix
         from sklearn.model selection import train test split
         from sklearn.neural network import MLPClassifier
         import sys
         from tensorflow import keras
         from tensorflow.keras.callbacks import LearningRateScheduler
         from tensorflow.keras.layers import Conv2D, BatchNormalization, Dense, Dropout, Flatten, MaxPool2D
         from tensorflow.keras.preprocessing.image import ImageDataGenerator
         from timeit import default timer as timer
         from utils import *
         mpl.rcParams.update({'font.size': 13})
         print('Python version', sys.version.split()[0])
```

Python version 3.8.7

# Import and preprocess data

The data files train.csv and test.csv contain grey-scale images of handwritten digits, from zero to nine. The training and test sets have 42000 and 28000 images, respectively. Each image has 28 pixels in height and 28 pixels in width, with 784 pixels in total. Each pixel-value is an integer between 0 and 255, inclusive. The first column of the training set is the digit label.

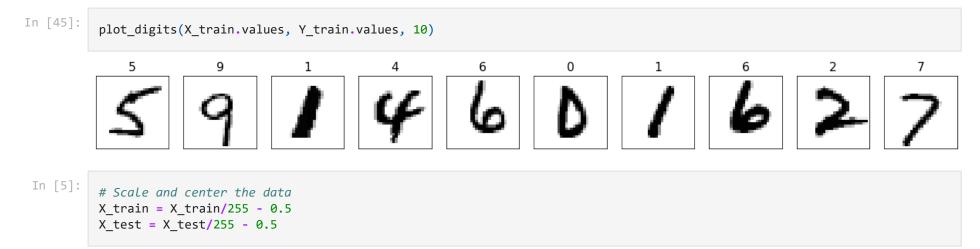
```
In [2]: # Import training and test sets
          folder = './'
          # folder = '/kaggle/input/digit-recognizer/'
          train = pd.read csv(folder + 'train.csv')
          X train = train.iloc[:,1:]
          Y train = train.iloc[:,0]
          X test = pd.read csv(folder + 'test.csv')
          n px = int(X train.shape[1]**0.5)
          print('Training set shape:', X_train.shape)
          print('Test set shape:', X_test.shape)
          print('N. of pixels', n px)
         Training set shape: (42000, 784)
         Test set shape: (28000, 784)
         N. of pixels 28
In [44]:
          def plot digits(X, Y, n digits):
              '''Plots the specified number of digits randomly selected from the data set and corresponding labels.'''
              # Select a random sample of digits
              idxs = np.random.randint(0, X.shape[0], n digits)
              X s = X[idxs]
              Y s = Y[idxs]
              # Number of rows and columns
              n cols = 10
              n rows = math.ceil(n digits/n cols)
              fig, axs = plt.subplots(n rows, n cols, figsize=(20,2.5*n rows), squeeze=False)
              i digit = 0
              for i row in range(n rows):
                  for i col in range(n cols):
                      ax = axs[i_row,i_col]
                      if i_digit >= n_digits:
                          ax.axis('off')
                          continue
                      # Reshape the pixels vector into a matrix
```

```
px = X_s[i_digit].reshape((n_px,-1))

# Plot the digit
ax.imshow(px, cmap='gray_r')
ax.set_xticks([])
ax.set_yticks([])
ax.set_title(Y_s[i_digit])

i_digit += 1
```

The following images show some handwritten digits and the corresponding labels.



We will split the current training set into a new training set and a validation set to assess the machine learning models' performance. We will use the train\_test\_split function of the Scikit-learn library. By dividing the initial data into three sets, we reduce the data available for training the model. Cross-validation) procedures can be used to mitigate this. Since a large data set is used in this project, and to keep a reasonable execution time of this notebook, this procedures won't be applied here.

```
In [6]:
# Split original training set into training and validation sets
X_train_o = X_train.copy()
Y_train_o = Y_train.copy()
X_train, X_val, Y_train, Y_val = train_test_split(X_train, Y_train, test_size=0.2, random_state=0)
```

### Model selection

We will try different machine learning models. Each algorithm has its strengths and weaknesses, and so we need to select the algorithm that works best with our specific data. We will compare the performance of the following algorithms:

- Logistic regression
- Fully connected feedforward neural network
- Convolutional neural network

The evaluation metric used to compare the different models is the categorization accuracy, i.e., the proportion of test images correctly classified.

#### Logistic regression

The logistic regression classifier is easy to implement and understand, so it is a natural first choice. We will use the LogisticRegression class of the Scikit-learn library.

```
In [7]:
         def fit_sklearn_model(model, X_train, Y_train, X_val=None, Y_val=None, display_info=True):
             Fits a Scikit-learn model to the training set and calculates its accuracy and training time.
             # Fit the model and calculate the training time
             start = timer()
             model.fit(X train, Y train)
             time = timer() - start
             # Calculate the accuracies
             train accuracy = accuracy score(Y train, model.predict(X train))
             if X val is not None:
                 val_accuracy = accuracy_score(Y_val, model.predict(X_val))
             else:
                 val accuracy = None
             # Display information of the training process
             if display info:
                 print('Train accuracy: {:.2%}'.format(train_accuracy), end='')
                 if X val is not None:
                     print('; Validation accuracy: {:.2%}'.format(val accuracy), end='')
                 print('; Training time: {:.1f}s'.format(time))
```

```
return train_accuracy, val_accuracy, time
```

```
In [8]: # Create the Logistic regression classifier
    create_arg = {'random_state':0,'solver':'newton-cg'}
    model = LogisticRegression(**create_arg)

# Fit the model
    fit_arg = {'X_train':X_train,'Y_train':Y_train,'X_val':X_val,'Y_val':Y_val}

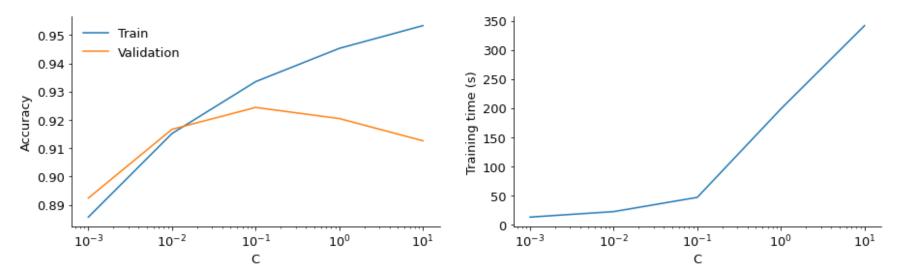
fit_sklearn_model(model, **fit_arg);
```

Train accuracy: 94.53%; Validation accuracy: 92.05%; Training time: 200.3s

The model's accuracy on the validation set is higher than that on the training set. The model might be overfitting the data, so we will try different normalisation levels to improve this. We will use five values, spaced evenly on a log scale between  $10^{-3}$  and 10.

```
In [9]:
         def compare_model_param(create_fun, create_arg, fit_fun, fit_arg, tuning_param, xscale='linear', xticks_labels=None):
             ''' Compares the models built with the specified hyperparameters.'''
             # Create and fit the models
             train accuracies = []
             val accuracies = []
             times = []
             for val in tuning param[1]:
                 model = create fun(**create arg, **{tuning param[0]:val})
                 outputs = fit_fun(model, **fit_arg)
                 train accuracies.append(get last elem(outputs[0]))
                 val accuracies.append(get last elem(outputs[1]))
                 times.append(outputs[2])
             # Plot the accuracies and training times
             if xticks labels is None:
                 x = tuning param[1]
             else:
                 x = range(len(tuning param[1]))
             fig = plt.figure(figsize=(15, 4))
```

```
plt.subplot(1, 2, 1)
              plt.plot(x, train accuracies)
              plt.plot(x, val accuracies)
              plt.xscale(xscale)
              plt.xlabel(tuning param[0])
              plt.xticks(x, xticks labels)
              plt.ylabel('Accuracy')
              plt.legend(['Train', 'Validation'], frameon=False)
              set spines()
              plt.subplot(1, 2, 2)
              plt.plot(x, times)
              plt.xscale(xscale)
              plt.xlabel(tuning param[0])
              plt.xticks(x, xticks labels)
              plt.ylabel('Training time (s)')
              set spines()
In [10]:
          C = np.logspace(-3, 1, 5)
          compare model param(LogisticRegression, create arg, fit sklearn model, fit arg, ('C',C), 'log')
         Train accuracy: 88.57%; Validation accuracy: 89.24%; Training time: 13.1s
         Train accuracy: 91.52%; Validation accuracy: 91.67%; Training time: 22.5s
         Train accuracy: 93.35%; Validation accuracy: 92.44%; Training time: 46.9s
         Train accuracy: 94.53%; Validation accuracy: 92.05%; Training time: 198.7s
         C:\PortablePrograms\WinPython\3 8\python-3.8.7.amd64\lib\site-packages\sklearn\utils\optimize.py:202: ConvergenceWarning:
         newton-cg failed to converge. Increase the number of iterations.
           warnings.warn("newton-cg failed to converge. Increase the "
         Train accuracy: 95.33%; Validation accuracy: 91.26%; Training time: 341.5s
```



A value of 0.1 led to the most accurate model. The LogisticRegressionCV class of the Scikit-learn library can also be used to select the best C parameter using a K-Folds cross-validator. This approach might lead to slightly better results but might be significantly more computationally expensive.

Now let's calculate the predictions for the test set.

```
In [11]:
# Create and fit the model using the optimised hyperparameter and the original training set
create_arg['C'] = 0.1

model = LogisticRegression(**create_arg)
fit_sklearn_model(model, X_train_o, Y_train_o);

# Predict class labels for test set and save them to file
def save_submission(Y_pred, filename):
    submission = pd.Series(Y_pred, index=range(1, len(Y_pred) + 1), name='Label')
    submission.to_csv(filename, index_label='ImageId')

save_submission(model.predict(X_test), 'submission_1.csv')
```

Train accuracy: 93.34%; Training time: 61.5s

The test set predictions were submitted to Kaggle, and an accuracy of 92.10% was obtained.

#### Fully connected feedforward neural network

Neural networks (NN) can model more complex structures in the data. However, there is generally a trade-off between interpretability and predictive power. The fully connected feedforward neural network was the first and most straightforward type. We will use the MLPClassifier class of the Scikit-learn library. We will start with one hidden layer with 100 units.

```
In [12]: # Create the multi-layer perceptron classifier
    create_arg = {'solver':'sgd','random_state':0}
    model = MLPClassifier(**create_arg)

# Fit the model
    fit_sklearn_model(model, **fit_arg);

Train accuracy: 97.19%; Validation accuracy: 95.79%; Training time: 148.8s

C:\PortablePrograms\WinPython\3_8\python-3.8.7.amd64\lib\site-packages\sklearn\neural_network\_multilayer_perceptron.py:6
14: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't converged yet.
    warnings.warn(
```

Using a NN with default hyperparameters instead of the initial logistic regression classifier increased the accuracy on the validation set from 92.05% to 95.79% while being significantly faster. Now let's tune the initial value of the learning rate and keep that value constant during training.

```
In [13]:
          def compare mlp loss(create fun, create arg, fit fun, fit arg, tuning param, legend=None):
              # Set default legend
              if legend is None:
                  legend = [str(val) for val in tuning param[1]]
              # Create and fit the models
              for i, val in enumerate(tuning param[1]):
                  model = create fun(**create arg, **{tuning param[0]:val})
                  time = fit fun(model, **fit arg, display info=False)[-1]
                  plt.plot(range(1, len(model.loss curve ) + 1), model.loss curve )
                  print('Value: {}; Training time: {:.1f}s'.format(legend[i], time))
              plt.xlabel('Iteration')
              plt.ylabel('Loss')
              plt.legend(legend, frameon=False)
              plt.title(tuning param[0])
              set spines()
```

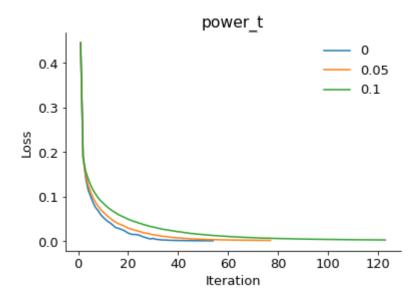
```
In [14]:
          # Tune the initial learning rate
          compare mlp loss(MLPClassifier, create arg, fit sklearn model, fit arg, ('learning rate init', [0.001,0.01,0.1]))
         C:\PortablePrograms\WinPython\3 8\python-3.8.7.amd64\lib\site-packages\sklearn\neural network\ multilayer perceptron.py:6
         14: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't converged yet.
           warnings.warn(
         Value: 0.001; Training time: 148.7s
         Value: 0.01; Training time: 139.9s
         Value: 0.1; Training time: 39.4s
                               learning rate init
            1.75 -
                                                          0.001
            1.50
                                                          0.01
                                                          0.1
            1.25
            1.00
         0.75
            0.50
            0.25
            0.00
                                       100
                                                  150
                   0
                             50
                                                             200
                                     Iteration
```

A value of 0.1 leads to the lowest loss and number of iterations. We can gradually decrease the learning rate during training to obtain more accurate models or a smoother variation of the loss function. We will use an inverse scaling exponent and tune the exponent value.

```
In [15]: # Tune the Learning rate schedule
    create_arg['learning_rate_init'] = 0.1
    create_arg['learning_rate'] = 'invscaling'
    vals = [0,0.05,0.1]

    compare_mlp_loss(MLPClassifier, create_arg, fit_sklearn_model, fit_arg, ('power_t',vals))

Value: 0; Training time: 39.5s
    Value: 0.05; Training time: 56.1s
    Value: 0.1; Training time: 90.0s
```

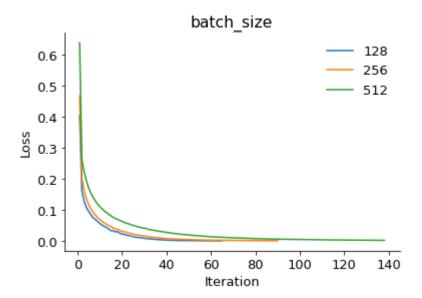


A value of 0.05 led to the best compromise between accuracy and smooth variation. Now let's tune the batch size, i.e., the number of samples per gradient update.

```
In [16]:
# Tune the batch size
    create_arg['power_t'] = 0.05

compare_mlp_loss(MLPClassifier, create_arg, fit_sklearn_model, fit_arg, ('batch_size',[128,256,512]))

Value: 128; Training time: 48.2s
    Value: 256; Training time: 62.3s
    Value: 512; Training time: 83.2s
```

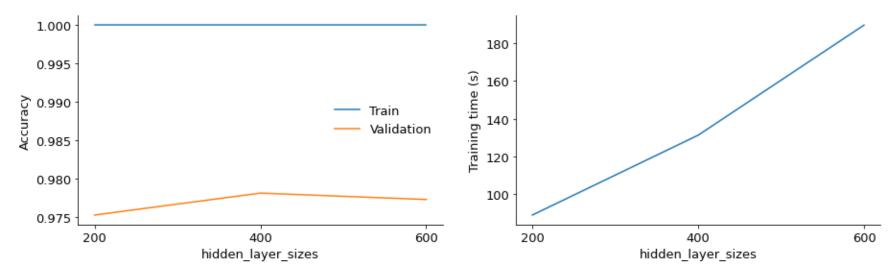


A value of 256 led to the best compromise between loss and smooth variation. Now let's tune the number of hidden layer's units.

```
In [17]:
    create_arg['batch_size'] = 256

    vals = [(200,),(400,),(600,)]
    compare_model_param(MLPClassifier, create_arg, fit_sklearn_model, fit_arg, ('hidden_layer_sizes',vals))

Train accuracy: 100.00%; Validation accuracy: 97.52%; Training time: 89.1s
    Train accuracy: 100.00%; Validation accuracy: 97.81%; Training time: 131.3s
    Train accuracy: 100.00%; Validation accuracy: 97.73%; Training time: 189.5s
```



Using 400 neurons lead to the highest accuracy on the validation set. Let's now try a NN with two layers.

```
In [18]:
           vals = [(400,200),(400,300),(400,400)]
           compare model param(MLPClassifier, create arg, fit sklearn model, fit arg, ('hidden layer sizes',vals),
                                xticks labels=[str(val) for val in vals])
          Train accuracy: 100.00%; Validation accuracy: 97.61%; Training time: 111.2s
          Train accuracy: 100.00%; Validation accuracy: 97.80%; Training time: 105.9s
          Train accuracy: 100.00%; Validation accuracy: 97.24%; Training time: 147.5s
             1.000
             0.995
                                                                                140
                                                                             Training time (s)
            0.990
          4ccuracy
                                                                                130
                                                              Train
             0.985
                                                              Validation
                                                                                120
             0.980
             0.975
                                                                                110
                (400, 200)
                                         (400, 300)
                                                                                                           (400, 300)
                                                                                                                                   (400, 400)
                                                                  (400, 400)
                                                                                  (400, 200)
                                     hidden layer sizes
                                                                                                      hidden layer sizes
```

The model with a second layer with 300 neurons led to the best results. The model's accuracy on the validation set is significantly higher

than that on the training set. The model might be overfitting the data, so we will try different normalisation values to improve this.

```
In [19]:
           create arg['hidden layer sizes'] = (400,300)
           compare model param(MLPClassifier, create arg, fit sklearn model, fit arg, ('alpha',np.logspace(-4, 0, 3)), 'log')
          Train accuracy: 100.00%; Validation accuracy: 97.80%; Training time: 105.1s
          Train accuracy: 100.00%; Validation accuracy: 97.86%; Training time: 144.9s
          Train accuracy: 97.49%; Validation accuracy: 96.30%; Training time: 338.7s
             1.00
                                                                 Train
                                                                 Validation
                                                                                    300
             0.99
                                                                                 Training time (s)
                                                                                    250
          Accuracy
              0.98
                                                                                    200
             0.97
                                                                                    150
                                                                                    100
                                             10^{-2}
                                                                        10°
                                                                                                                   10^{-2}
                                                                                                                                              10<sup>0</sup>
                   10^{-4}
                                                                                         10^{-4}
                                             alpha
                                                                                                                   alpha
```

Adding normalization didn't improve the accuracy of the model, and so we will assume that the model is not overfitting the data. We could have used the classes GridSearchCV or RandomizedSearchCV of the Scikit-learn library to tune the hyperparameters. These approaches might lead to more optimised hyperparameters but might be more computationally expensive. Also, the approach we followed allows a better understanding of how the different parameters influence our model's performance.

Now let's calculate the predictions for the test set.

```
# Create and fit the model using the optimised hyperparameters and the original training set
model = MLPClassifier(**create_arg)
fit_sklearn_model(model, X_train_o, Y_train_o);

# Predict class labels for test set and save them to file
save_submission(model.predict(X_test), 'submission_2.csv')
```

Train accuracy: 100.00%; Training time: 126.0s

The test set predictions were submitted to Kaggle, and an accuracy of 97.85% was obtained.

#### Convolutional neural network

Convolutional Neural Networks (CNN) have demonstrated superior results in image classification. A typical CNN architecture consists of pairs of convolution and pooling layers, followed by dense layers and a final softmax layer.

We will use Keras library to develop our model. We need to reshape the data sets so that the pixels of each image are stored as a (28,28,1) array. Also, we need to convert the class vectors to binary class matrixes using the function to\_categorical.

```
In [21]: # Reshape the data sets
X_train_r = X_train.values.reshape(-1,n_px,n_px,1)
X_val_r = X_val.values.reshape(-1,n_px,n_px,1)
X_test_r = X_test.values.reshape(-1,n_px,n_px,1)

# Convert the class vectors to binary class matrixes (one-hot encoding)
Y_train_oh = keras.utils.to_categorical(Y_train, num_classes=10)
Y_val_oh = keras.utils.to_categorical(Y_val, num_classes=10)
```

There is an infinity number of possible CNN architectures, and researchers have spent vasts amounts of time developing different architectures that are optimised for various applications. Hence, as a first step, we should generally search the current literature for an architecture applied in an application similar to ours. We will start with an architecture similar to the LeNet-5, which has been developed for handwriting recognition (LeCun et al., 1998).

The function Sequential is used to initialize our model. The convolution, pooling, dropout and dense layers are defined using the functions Conv2D, MaxPool2D, Dropout and Dense. The Adam optimizer is used for training the model.

```
def create_keras_model(filters=(8,16), kernel_size=5, units=256, dropout_rate=(0,0,0), learning_rate=0.001):
    '''Creates a Keras model using the specified hyperparameters.'''

# Initialize the model
    model = keras.Sequential()

# Add Layers
    model.add(Conv2D(filters=filters[0], kernel_size=kernel_size, padding='same', activation='relu', input_shape=(n_px,n_model.add(MaxPool2D()))
    model.add(Dropout(dropout_rate[0]))
    model.add(Conv2D(filters=filters[1], kernel_size=kernel_size, padding='same', activation='relu'))
    model.add(MaxPool2D())
```

```
model.add(Dropout(dropout_rate[1]))
model.add(Flatten())
model.add(Dense(units, activation='relu'))
model.add(Dropout(dropout_rate[2]))
model.add(Dense(10, activation='softmax'))

# Compile the model
model.compile(keras.optimizers.Adam(learning_rate), 'categorical_crossentropy', ['accuracy'])
return model
```

The following function is used to fit Keras models.

```
In [23]:
          def fit_keras_model(model, X_train, Y_train, datagen=None, X_val=None, Y_val=None, batch_size=64, epochs=30, callbacks=None
                              display info=True):
              Fits a Keras model to the training set and calculates its accuracy and training time.
              # Fit the model and calculate the training time
              start = timer()
              if X val is None:
                  data val = None
              else:
                  data val = (X val,Y val)
              if datagen is None:
                  history = model.fit(X_train, Y_train, batch_size=batch_size, epochs=epochs, callbacks=callbacks,
                                      validation data=data val, verbose=0).history
              else:
                  history = model.fit(datagen.flow(X train, Y train, batch size=batch size), batch size=batch size, epochs=epochs,
                                       callbacks=callbacks, validation data=data val, verbose=0).history
              time = timer() - start
              train_accuracy = history['accuracy']
              if X val is None:
                  val accuracy = None
              else:
                  val accuracy = history['val accuracy']
              # Display information of the training process
              if display info:
                  print('Train accuracy: {:.2%}'.format(train accuracy[-1]), end='')
```

```
if X_val is not None:
    print('; Validation accuracy: {:.2%}'.format(val_accuracy[-1]), end='')
print('; Training time: {:.1f}s'.format(time))

return train_accuracy, val_accuracy, time
```

We will start by tuning the initial value of the learning rate and keep that value constant during training.

```
In [25]:
# Tune the initial Learning rate
fit_arg = {'X_train':X_train_r,'Y_train':Y_train_oh,'X_val':X_val_r,'Y_val':Y_val_oh}

legend = []

for val in [1E-4,1E-3,1E-2]:

    model = create_keras_model(learning_rate=val)

    train_accuracy = fit_keras_model(model, **fit_arg)[0]

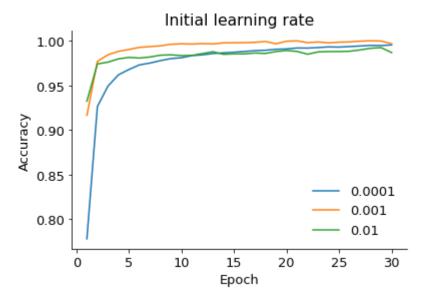
    plt.plot(range(1, len(train_accuracy) + 1), train_accuracy)

    legend.append(str(val))

plt.xlabel('Epoch')
    plt.ylabel('Accuracy')
    plt.legend(legend, frameon=False)
    plt.title('Initial learning rate')
    set_spines()

Train accuracy: 99.53%; Validation accuracy: 98.61%; Training time: 279.2s
```

Train accuracy: 99.53%; Validation accuracy: 98.61%; Training time: 279.2s Train accuracy: 99.66%; Validation accuracy: 98.44%; Training time: 282.6s Train accuracy: 98.67%; Validation accuracy: 97.71%; Training time: 275.4s



Using a value of  $10^{-4}$  leads to a more smooth variation but needs many more epochs to converge to a value near the optimum. We will use a value of  $10^{-3}$  and tune the learning rate's decay rate to get a smoother variation. Also, using 20 epochs to train the model seems enough.

```
In [26]: create_arg = {'learning_rate':0.001}
fit_arg['epochs'] = 20

In [27]: def compare_keras_fit_param(create_fun, create_arg, fit_fun, fit_arg, tuning_param, legend=None):
    '''Compares the models built with the specified hyperparameters.'''
    if legend == None:
        legend = [str(val) for val in tuning_param[1]]
        for val in tuning_param[1]:
            model = create_fun(**create_arg)
            train_accuracy = fit_fun(model, **fit_arg, **{tuning_param[0]:val})[0]
            plt.plot(range(1, len(train_accuracy) + 1), train_accuracy)
            plt.xlabel('Epoch')
            plt.ylabel('Accuracy')
            plt.legend(legend, frameon=False)
```

A decay rate of 0.95 leads to the best results. Now let's tune the batch size.

Epoch

10

5

Accuracy

0.96

0.94

0.92

```
In [29]: # Tune the batch size
    fit_arg['callbacks'] = schedulers[1]

compare_keras_fit_param(create_keras_model, create_arg, fit_keras_model, fit_arg, ('batch_size',[32,64,128]))

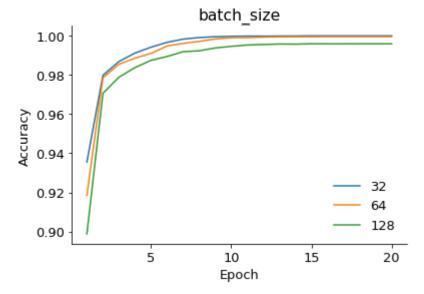
Train accuracy: 100.00%; Validation accuracy: 98.96%; Training time: 224.5s
Train accuracy: 99.95%; Validation accuracy: 98.94%; Training time: 185.4s
Train accuracy: 99.59%; Validation accuracy: 98.65%; Training time: 170.3s
```

0.95

20

0.9

15

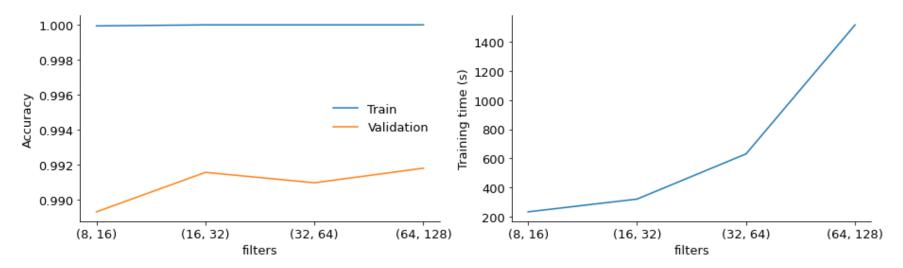


A batch size of 32 yields the best results.

```
In [55]:
fit_arg['batch_size'] = 32
```

We will now tune the following hyperparameters used for creating the models:

- Number of filters of the convolution layers;
- Number of units of the first dense layer;
- Kernel size of the convolution layers;
- Rates of the dropout layers.



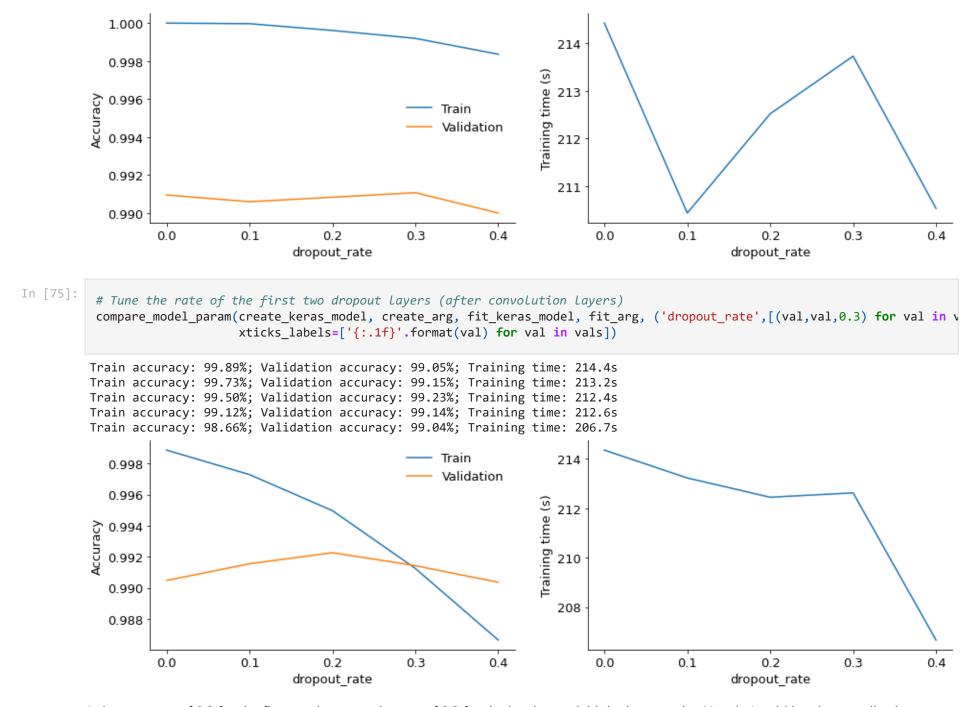
Using 16 and 32 filters leads to the best results.

```
In [62]:
           create_arg['filters'] = (16,32)
In [70]:
           # Tune the number of units of the first dense layer
           compare_model_param(create_keras_model, create_arg, fit_keras_model, fit_arg, ('units',[128,256,512]))
          Train accuracy: 100.00%; Validation accuracy: 99.00%; Training time: 291.7s
          Train accuracy: 100.00%; Validation accuracy: 99.10%; Training time: 323.8s
          Train accuracy: 100.00%; Validation accuracy: 99.11%; Training time: 373.5s
            1.000
                                                                                360
             0.998
                                                                             Training time (s)
          Accuracy
             0.996
                                                                                340
                                                              Train
                                                              Validation
             0.994
                                                                                320
             0.992
                                                                                300
             0.990
                                    256
                                                                     512
                                                                                     128
                                                                                                     256
                                                                                                                                      512
                    128
                                            units
                                                                                                             units
```

Using 256 units yields the best results without compromising the training time.

```
In [71]:
           create arg['units'] = 256
In [72]:
           # Tune the kernel size of the convolution layers
           compare model param(create keras model, create_arg, fit_keras_model, fit_arg, ('kernel_size',[3,5,7]))
          Train accuracy: 100.00%; Validation accuracy: 99.15%; Training time: 208.9s
          Train accuracy: 100.00%; Validation accuracy: 99.17%; Training time: 324.1s
          Train accuracy: 100.00%; Validation accuracy: 99.13%; Training time: 498.8s
                                                                                 500
             1.000
                                                                                 450
             0.998
                                                                              raining time (s)
                                                                                 400
          Accuracy
                                                               Train
             0.996
                                                                                 350
                                                               Validation
             0.994
                                                                                 300
                                                                                 250
             0.992
                                                                                 200
                     3
                                              5
                                                                                                                5
                                                                                                                                         7
                                         kernel size
                                                                                                           kernel size
```

A kernel of size 3 seems a good compromise between accuracy and efficiency.

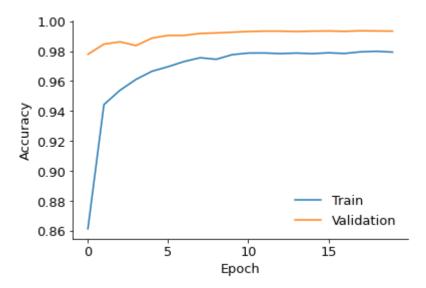


A dropout rate of 0.2 for the first two layers, and a rate of 0.3 for the last layer, yield the best results. Now let's add batch normalization layers and data augmentation to improve the performance of our model. The Keras class ImageDataGenerator is used for generating

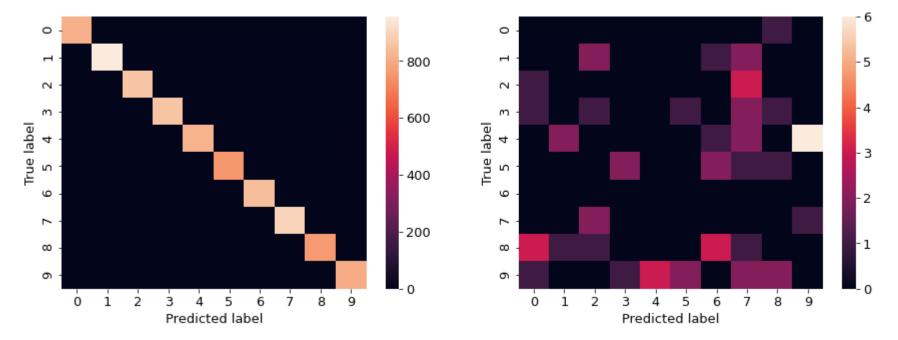
image data, and the parameters used are described in this notebook done by Chris Deotte.

```
In [76]:
          # Create the model
          model = keras.Sequential()
          model.add(Conv2D(filters=16, kernel size=3, padding='same', activation='relu', input shape=(n px,n px,1)))
          model.add(BatchNormalization())
          model.add(MaxPool2D())
          model.add(Dropout(0.2))
          model.add(Conv2D(filters=32, kernel size=3, padding='same', activation='relu'))
          model.add(BatchNormalization())
          model.add(MaxPool2D())
          model.add(Dropout(0.2))
          model.add(Flatten())
          model.add(Dense(256, activation='relu'))
          model.add(BatchNormalization())
          model.add(Dropout(0.3))
          model.add(Dense(10, activation='softmax'))
          model.compile(keras.optimizers.Adam(0.001), 'categorical crossentropy', ['accuracy'])
          # Create image data generator
          datagen = ImageDataGenerator(rotation range=10, zoom range=0.1, width shift range=0.1, height shift range=0.1)
          # Fit the model using the optimised hyperparameters
          train accuracy, val accuracy, = fit keras model(model, X train r, Y train oh, datagen, X val r, Y val oh, 32, 20,
                                                             schedulers[1])
          plt.plot(train accuracy)
          plt.plot(val accuracy)
          plt.xlabel('Epoch')
          plt.ylabel('Accuracy')
          plt.legend(['Train', 'Validation'], frameon=False)
          set spines()
```

Train accuracy: 97.93%; Validation accuracy: 99.33%; Training time: 443.5s



Now let's look at the wrongly predicted digits. The confusion matrix is shown on the left subplot of the figure below. Because the number of digits predicted correctly is much higher than those mislabeled, we mainly see the matrix's diagonal. Its colour does not vary much because we have a uniform distribution of digits' types. The right subplot shows the confusion matrix with its diagonal filled with zeros to highlight the wrong predictions.



The first row of the following matrix is the predicted label, and the second row is the sum of the elements of each column of the confusion matrix. E.g., 13 digits were wrongly predicted to be 7.

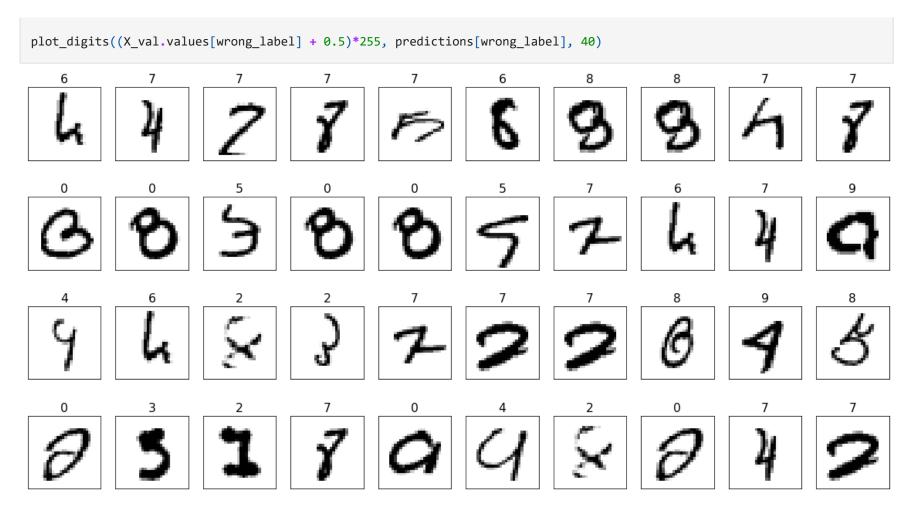
```
In [117...
    print(np.vstack((np.arange(10),cm.sum(axis=0))))

[[ 0 1 2 3 4 5 6 7 8 9]
    [ 6 3 6 3 3 3 7 13 5 7]]
```

The first row of the following matrix is the true label, and the second row is the sum of the elements of each row of the confusion matrix. E.g., the digits 4 and 9 have 11 wrong predictions each.

Now let's plot 40 digits, randomly selected, that have wrong predictions. The title of each plot is the predicted label. We can observe several of the cases mentioned before. E.g., several times, the digit 2 was wrongly predicted to be 7.

```
In [77]:
    predictions = model.predict(X_val_r).argmax(axis=1)
    wrong_label = Y_val != predictions
```



Let's calculate the predictions for the test set.

Train accuracy: 98.18%; Training time: 673.0s

The test set predictions were submitted to Kaggle, and an accuracy of 99.25% was obtained.

## Conclusions and further developments

The following table shows the accuracies and training times of the different types of models. The logistic regression is clearly not a good choice for this type of applications. Using a CNN leads to the best results but also requires much more time to train. It would be interesting to tune a smaller CNN and see if we could obtain an accuracy and efficiency better then those of the fully connected NR. The accuracy of the CNN can be further improved by using ensemble methods and collecting more data. For the latter, we recommend paying particular attention to the digits 4, 7 and 9.

Model	Accuracy (%)	Training time (s)
Logistic regression	92.10	61.5
Fully connected neural network	97.85	126.0
Convolutional neural network	99.25	673.0

# Acknowledgements

These excellent notebooks inspired some of the code used in this notebook:

- MNIST Logistic Regression Baseline by Bojan Tunguz
- How to choose CNN Architecture MNIST by Chris Deotte
- How to score 97%, 98%, 99%, and 100% by Chris Deotte