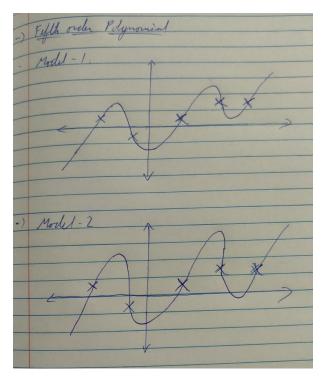
Assignment- 4

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Exercise: Draw some points on a 2D plane to which a fifth-order polynomial can fit. Show one candidate fifth order polynomial that will fit these points. Now show an alternate fifth order polynomial that will fit the same points. Comment on the relative coefficients of the two polynomials (how can you do this from a sketch?).



We can see that both the models fit the data points and are both fifth order polynomials but model 1 has smaller slopes than model 2 which implies that the coefficients of the first model will be smaller than the second model.

We can also say that the first model is a better fit as it is less prone to overfitting whereas the second model is prone to overfitting the dataset

Exercise: Contrast this exercise with the motivating example from lecture I on polynomial regression.

In this exercise we can see even by keeping the DOF's constant the fit and generalization of the model depends on its weights- how large they may be. Larger weights may overfit the dataset

while smaller weights may be more generalized. Whereas in the polynomial regression we increases the DOF's step by step in order to minimize the possibility of the model to overfit the data.

Exercise: Why are we only proposing augmentation of the training dataset?

Augmentation produces synthetic data that may have a bias imparted by the creator of the synthetic data. While training this may be acceptable as we are trying to train our model with as many possibilities possible but while we validate the model we want to validate it with proper data and not make any assumptions on our data set.

Exercise: Does a bottleneck layer require nonlinearities? Justify your answer

The bottleneck layer doesn't need Non linearities as the main function of this layer is to reduce the number of DOF's in the system and the data from layer A will be transferred to layer B through this bottle neck layer. For this purpose a linear function can be used in the bottle neck layer as it will work fine for this purpose.

Exercise: Suppose we knew that model M1 with N1 DOFs fit the first mode of operation, and model M2 with N2 DOFs fit the second mode of operation. Describe how these three solutions {M,N}, {M1,N1} and {M2,N2} relate to each other in terms of suitability to solving the problem, lower number of DOFs, etc. Consider a practical case: suppose our dataset expresses a relationship that has (for some inputs) a fourth order polynomial nonlinear relationship, and (for other inputs) a linear relationship. How would {M,N}, {M1,N1}, {M2,N2} look for this case?

M2 and M1 working together will give us a better output than a single model M as the DOF required by M (N) will be larger than N1 and N2 which in turn will be more prone to overfitting than models M1 and M2.

In the practical case we can say the minimum DOF needed for M1 to fit a 4^{th} order polynomial is 5-4 weights and a bias and the minimum DOF for M2 to fit a linear polynomial is 2-1 weight and a bias. Whereas a model M to fit to fit both the 4^{th} order polynomial and the linear polynomial will be 6 or greater and also depends on the interaction of these two modes of operation with each other and may requires many DOF's to truly represent their relationship

and properly fit both modes. Hence can be prone to overfitting as we increase the DOF of model M.

Question- Dropout

1. Network structure

1-)	Network Structure. 3
-	28 317 -> 31
	maye 32 filter 69 filter
-	TO SEC FICE
	Max Pool Fletten loyer 4 Output loyer 33,856 128 layer
	Neuron Neuron 10 Neuron

- CNN Dropout

```
# TensorFlow and tf.keras
import tensorflow as tf
from tensorflow import keras
from tensorflow.keras.layers import Conv2D, MaxPooling2D, Dropout, Flatten, Dense
# Commonly used modules
import numpy as np
import os
import sys
# Images, plots, display, and visualization
import matplotlib.pyplot as plt
import pandas as pd
import seaborn as sns
import cv2
import IPython
from six.moves import urllib
print(tf. version )
 □→ 2.4.1
(train images, train labels), (test images, test labels) = keras.datasets.mnist.load data()
# reshape images to specify that it's a single channel
print(test_labels.shape)
train images = train images.reshape(train images.shape[0], 28, 28, 1)
test_images = test_images.reshape(test_images.shape[0], 28, 28, 1)
     Downloading data from <a href="https://storage.googleapis.com/tensorflow/tf-keras-datasets/mnist.npz">https://storage.googleapis.com/tensorflow/tf-keras-datasets/mnist.npz</a>
     (10000,)
```

```
det preprocess images(imgs): # should work tor both a single image and multiple images
    sample img = imgs if len(imgs.shape) == 2 else imgs[0]
    assert sample_img.shape in [(28, 28, 1), (28, 28)], sample_img.shape # make sure images are 28x28 and single-channel (gra
    return imgs / 255.0
train_images = preprocess_images(train_images)
test images = preprocess images(test images)
plt.figure(figsize=(10,2))
for i in range(10,15):
    plt.subplot(1,5,i-9)
    plt.xticks([])
    plt.yticks([])
    plt.grid(False)
    plt.imshow(train images[i].reshape(28, 28), cmap=plt.cm.binary)
    plt.xlabel(train labels[i])
```

2. Annotate the sizes of the mathematical objects within; inputs, outputs, filters, etc.

The input to the neural network are 28 X 28 images- the train set has 60000 images and test set has 10000 images.

The output vector is a 10 element one hot vector that maps the output to numbers ranging from 0-9

The model-

- 1 layer- 32 convolution filters of size 3 x 3
- 2 layer- 64 convolution filters of size 3 x 3 x 32
- 3 layer- max-pool filter of size 2 x 2

- 4 layer- fully conected layer with 128 neurons connnected with max pool layer
- Output layer- 10 neurons fully connected with layer 4

The ouputs of each layer and the weights of fully connected layers-

- 1 layer-26 x 26 x 32
- 2 layer- 24 x 24 x 64
- 3 layer- 23 x 23 x 64
- 4 layer- 128 outputs from 33,856 inputs- which gives us 4,333,568 weights and 128 biases
 - Output layer- 10 outputs from 128 inputs- which gives us 1,280 weights and 10 biases

```
model = keras.Sequential()
# 32 convolution filters used each of size 3x3
model.add(Conv2D(32, kernel_size=(3, 3), activation='relu', input_shape=(28, 28, 1)))
# 64 convolution filters used each of size 3x3
model.add(Conv2D(64, (3, 3), activation='relu'))
# choose the best features via pooling
model.add(MaxPooling2D(pool size=(2, 2)))
# randomly turn neurons on and off to improve convergence
model.add(Dropout(0.25))
# flatten since too many dimensions, we only want a classification output
model.add(Flatten())
# fully connected to get all relevant data
model.add(Dense(128, activation='relu'))
# one more dropout
model.add(Dropout(0.5))
# output a softmax to squash the matrix into output probabilities
model.add(Dense(10, activation='softmax'))
```

3. How is dropout applied?

Dropout is applied to the previous layer where the arrgument represents the chance with which a neurons activation will be 0. eg- 0.5 means that there is a 50% chance of dropout for each neuron in the layer before

```
model.compile(optimizer=tf.optimizers.Adam(),
       loss='sparse categorical crossentropy',
       metrics=['accuracy'])
history = model.fit(train images, train labels, epochs=5)
  Epoch 1/5
  1875/1875 [=============== ] - 145s 77ms/step - loss: 0.3605 - accuracy: 0.8879
  Epoch 2/5
  Epoch 3/5
  Epoch 4/5
  Epoch 5/5
  print(test images.shape)
test loss, test acc = model.evaluate(test_images, test_labels)
print('Test accuracy:', test acc)
  (10000, 28, 28, 1)
  Test accuracy: 0.9902999997138977
```

4. What is the performance with Dropout enabled?

There is very high accurancy for both the training set and the test set while using drop out. The varience between both the training accuracy and the test accuracy is also quite low

Removing Dropout

```
model2 = keras.Sequential()
  # 32 convolution filters used each of size 3x3
  model2.add(Conv2D(32, kernel size=(3, 3), activation='relu', input shape=(28, 28, 1)))
  # 64 convolution filters used each of size 3x3
  model2.add(Conv2D(64, (3, 3), activation='relu'))
  # choose the best features via pooling
  model2.add(MaxPooling2D(pool size=(2, 2)))
  # flatten since too many dimensions, we only want a classification output
  model2.add(Flatten())
  # fully connected to get all relevant data
  model2.add(Dense(128, activation='relu'))
  # output a softmax to squash the matrix into output probabilities
  model2.add(Dense(10, activation='softmax'))
  model2.compile(optimizer=tf.optimizers.Adam(),
            loss='sparse categorical crossentropy',
            metrics=['accuracy'])
  history = model2.fit(train images, train labels, epochs=5)
      Epoch 1/5
      Epoch 2/5
      Epoch 3/5
      Epoch 4/5
      Epoch 5/5
      print(test images.shape)
  test loss, test acc = model2.evaluate(test images, test labels)
https://colab.research.google.com/drive/1VJza5Cr1RF6IBmvHppO-jJs6EK2gqXby#scrollTo=OYrk7FIh7ytz&printMode=true
```

5. Now disable it. What is the performance?

Disabling dropout we can notice the accuracy of the training set increased but there is a slighty larger gap in accuracy between the training set and the test set. This may imply a case of slight overfitting to the Training set.

Increasing the probability of Dropout

```
model3 = keras.Sequential()
# 32 convolution filters used each of size 3x3
model3.add(Conv2D(32, kernel size=(3, 3), activation='relu', input shape=(28, 28, 1)))
# 64 convolution filters used each of size 3x3
model3.add(Conv2D(64, (3, 3), activation='relu'))
# choose the best features via pooling
model3.add(MaxPooling2D(pool size=(2, 2)))
# randomly turn neurons on and off to improve convergence
model3.add(Dropout(0.5))
# flatten since too many dimensions, we only want a classification output
model3.add(Flatten())
# fully connected to get all relevant data
model3.add(Dense(128, activation='relu'))
# one more dropout
model3.add(Dropout(0.7))
# output a softmax to squash the matrix into output probabilities
model3.add(Dense(10, activation='softmax'))
model3.compile(optimizer=tf.optimizers.Adam(),
              loss='sparse categorical crossentropy',
```

```
metrics=['accuracy'])
```

```
history = model3.fit(train images, train labels, epochs=5)
 Epoch 1/5
 Epoch 2/5
 Epoch 3/5
 Epoch 4/5
 Epoch 5/5
 print(test images.shape)
test loss, test acc = model3.evaluate(test images, test labels)
print('Test accuracy:', test acc)
 (10000, 28, 28, 1)
 Test accuracy: 0.9898999929428101
```

Using a lower probability of Dropout

```
model4 = keras.Sequential()
# 32 convolution filters used each of size 3x3
model4.add(Conv2D(32, kernel_size=(3, 3), activation='relu', input_shape=(28, 28, 1)))
# 64 convolution filters used each of size 3x3
model4.add(Conv2D(64, (3, 3), activation='relu'))
# choose the best features via pooling
model4.add(MaxPooling2D(pool_size=(2, 2)))
# randomly turn neurons on and off to improve convergence
model4.add(Dropout(0.1))
# flatten since too many dimensions, we only want a classification output
```

```
model4.add(Flatten())
# fully connected to get all relevant data
model4.add(Dense(128, activation='relu'))
# one more dropout
model4.add(Dropout(0.1))
# output a softmax to squash the matrix into output probabilities
model4.add(Dense(10, activation='softmax'))
model4.compile(optimizer=tf.optimizers.Adam(),
        loss='sparse categorical crossentropy',
        metrics=['accuracy'])
history = model4.fit(train images, train labels, epochs=5)
   Epoch 1/5
   Epoch 2/5
   Epoch 3/5
   Epoch 4/5
   Epoch 5/5
   1875/1875 [================ ] - 153s 81ms/step - loss: 0.0077 - accuracy: 0.9973
print(test images.shape)
test loss, test acc = model4.evaluate(test images, test labels)
print('Test accuracy:', test acc)
   (10000, 28, 28, 1)
   Test accuracy: 0.9916999936103821
```

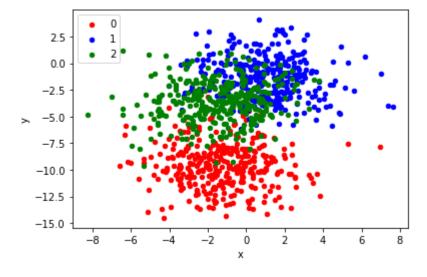
6 and 7

Altering the dropout in model 3 we can see that increasing drop out to a higher extent- Lowers the DOF of the model drastically while training as most of the neurons get turned off and the much of the useful data is lost. This leads to a lower accuracy in both the training and test set. We can also see a larger gap between the accuracy of the test and training set.

Altering the dropout in model 4 we can see that decreasing the dropout closer to 0 brings us closer to a model that doesnt have drop out but also gives us the benfits of generalization of dropout. The accuracy of the training set is close to the model with no dropout. The gap in the accuracy of the test set and the train set is also quite low; it is comparable to the first model.

We can see that there is no definate value of dropot that will give us the best model for our problem. This is why this is a hyperparameter of the model and has to be empirically choosen by us.

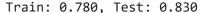
```
# TensorFlow and tf.keras
import tensorflow as tf
from tensorflow import keras
from tensorflow.keras.layers import Conv2D, MaxPooling2D, Dropout, Flatten, Dense
from sklearn.datasets import make blobs
from sklearn.metrics import accuracy score
from itertools import product
from numpy.linalg import norm
# Commonly used modules
import numpy as np
import os
import sys
# Images, plots, display, and visualization
import matplotlib.pyplot as plt
from pandas import DataFrame
# generate 2d classification dataset
X, y = make blobs(n samples=1000, centers=3, n features=2, cluster std=2, random state=2)
# scatter plot, dots colored by class value
df = DataFrame(dict(x=X[:,0], y=X[:,1], label=y))
colors = {0:'red', 1:'blue', 2:'green'}
fig, ax = plt.subplots()
grouped = df.groupby('label')
for key, group in grouped:
    group.plot(ax=ax, kind='scatter', x='x', y='y', label=key, color=colors[key])
plt.show()
```

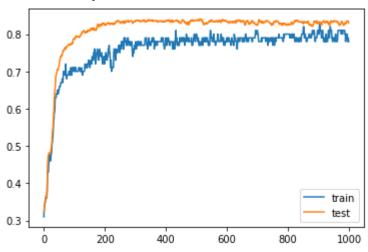


```
# one hot encode output variable
y1 = keras.utils.to_categorical(y)
```

```
# split into train and test
```

```
n_{\tau} = 100
trainX, testX = X[:n_train, :], X[n_train:, :]
trainy, testy = y1[:n_train], y1[n_train:]
print(trainX.shape, testX.shape)
     (100, 2) (900, 2)
# making the model
model = keras.Sequential()
model.add(Dense(25, input dim=2, activation='relu'))
model.add(Dense(3, activation='softmax'))
model.compile(loss='categorical_crossentropy', optimizer='adam', metrics=['accuracy'])
# fit model
history = model.fit(trainX, trainy, validation_data=(testX, testy), epochs=1000, verbose=0)
# evaluate the model
_, train_acc = model.evaluate(trainX, trainy, verbose=0)
_, test_acc = model.evaluate(testX, testy, verbose=0)
print('Train: %.3f, Test: %.3f' % (train_acc, test_acc))
# learning curves of model accuracy
plt.plot(history.history['accuracy'], label='train')
plt.plot(history.history['val_accuracy'], label='test')
plt.legend()
plt.show()
```



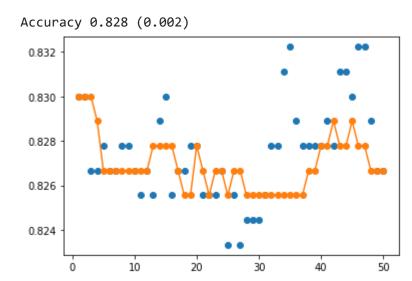


Horizontal Voting Ensemble-

```
# create directory for models
os.makedirs('model')
# fit model
n_epochs, n_save_after = 1000, 950
```

```
for i in range(n epochs):
 # fit model for a single epoch
  model.fit(trainX, trainy, epochs=1, verbose=0)
 # check if we should save the model
  if i >= n save after:
    model.save('models/model_' + str(i) + '.h5')
# load models from file
def load all models(n start, n end):
  all models = list()
  for epoch in range(n start, n end):
    # define filename for this ensemble
    filename = 'models/model ' + str(epoch) + '.h5'
    # load model from file
    model = keras.models.load model(filename)
    # add to list of members
    all models.append(model)
    print('>loaded %s' % filename)
  return all models
# make an ensemble prediction for multi-class classification
def ensemble predictions(members, testX):
 # make predictions
 yhats = [model.predict(testX) for model in members]
 yhats = np.array(yhats)
  # sum across ensemble members
  summed = np.sum(yhats, axis=0)
  # argmax across classes
  result = np.argmax(summed, axis=1)
  return result
# evaluate a specific number of members in an ensemble
def evaluate n members(members, n members, testX, testy):
  # select a subset of members
  subset = members[:n members]
  # make prediction
 yhat = ensemble_predictions(subset, testX)
  # calculate accuracy
  return accuracy_score(testy, yhat)
# load models in order
members = load all models(950, 1000)
print('Loaded %d models' % len(members))
# reverse loaded models so we build the ensemble with the last models first
members = list(reversed(members))
SHOW HIDDEN OUTPUT
```

```
# evaluate different numbers of ensembles on hold out set
single scores, ensemble scores = list(), list()
for i in range(1, len(members)+1):
 # evaluate model with i members
 ensemble_score = evaluate_n_members(members, i, testX, testy)
 # evaluate the i'th model standalone
 testy_enc = keras.utils.to_categorical(testy)
 _, single_score = members[i-1].evaluate(testX, testy_enc, verbose=0)
 # summarize this step
 print('> %d: single=%.3f, ensemble=%.3f' % (i, single_score, ensemble_score))
 ensemble scores.append(ensemble score)
 single_scores.append(single_score)
SHOW HIDDEN OUTPUT
# summarize average accuracy of a single final model
print('Accuracy %.3f (%.3f)' % (np.mean(single_scores), np.std(single_scores)))
# plot score vs number of ensemble members
x_axis = [i for i in range(1, len(members)+1)]
plt.plot(x_axis, single_scores, marker='o', linestyle='None')
plt.plot(x axis, ensemble scores, marker='o')
plt.show()
```

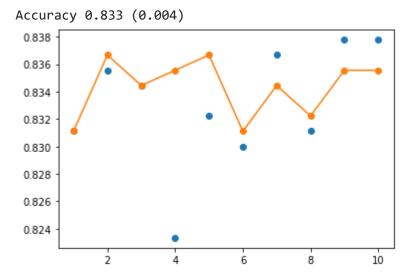


we can see that the ensemble accuracy is quite stable and has a lower varience than the single models.

Average Ensemble

```
# fit model on dataset
def fit model(trainX, trainy):
  trainy enc = keras.utils.to categorical(trainy)
  # define model
```

```
model = keras.Sequential()
 model.add(Dense(25, input dim=2, activation='relu'))
 model.add(Dense(3, activation='softmax'))
 model.compile(loss='categorical_crossentropy', optimizer='adam', metrics=['accuracy'])
 # fit model
 model.fit(trainX, trainy enc, epochs=500, verbose=0)
 return model
n members = 10
members = [fit model(trainX, trainy) for    in range(n members)]
# evaluate different numbers of ensembles on hold out set
single scores, ensemble scores = list(), list()
for i in range(1, len(members)+1):
 # evaluate model with i members
 ensemble score = evaluate n members(members, i, testX, testy)
 # evaluate the i'th model standalone
 testy enc = keras.utils.to categorical(testy)
 _, single_score = members[i-1].evaluate(testX, testy_enc, verbose=0)
 # summarize this step
 print('> %d: single=%.3f, ensemble=%.3f' % (i, single score, ensemble score))
 ensemble scores.append(ensemble score)
  single scores.append(single score)
     > 1: single=0.831, ensemble=0.831
     > 2: single=0.836, ensemble=0.837
     > 3: single=0.834, ensemble=0.834
     > 4: single=0.823, ensemble=0.836
     > 5: single=0.832, ensemble=0.837
     > 6: single=0.830, ensemble=0.831
     > 7: single=0.837, ensemble=0.834
     > 8: single=0.831, ensemble=0.832
     > 9: single=0.838, ensemble=0.836
     > 10: single=0.838, ensemble=0.836
# summarize average accuracy of a single final model
print('Accuracy %.3f (%.3f)' % (np.mean(single_scores), np.std(single_scores)))
# plot score vs number of ensemble members
x_axis = [i for i in range(1, len(members)+1)]
plt.plot(x axis, single scores, marker='o', linestyle='None')
plt.plot(x_axis, ensemble_scores, marker='o')
plt.show()
```



We can see using the average ensemble we get an accuracy that is neither too high nor too low there by reducing our models varience.

Grid Search Weighted Average Ensemble

```
# make an ensemble prediction for multi-class classification
def ensemble_predictions(members, weights, testX):
  # make predictions
  yhats = [model.predict(testX) for model in members]
  yhats = np.array(yhats)
  # weighted sum across ensemble members
  summed = np.tensordot(yhats, weights, axes=((0),(0)))
  # argmax across classes
  result = np.argmax(summed, axis=1)
  return result
# evaluate a specific number of members in an ensemble
def evaluate ensemble(members, weights, testX, testy):
  # make prediction
  yhat = ensemble predictions(members, weights, testX)
  # calculate accuracy
  return accuracy_score(testy, yhat)
# normalize a vector to have unit norm
def normalize(weights):
  # calculate l1 vector norm
  result = norm(weights, 1)
  # check for a vector of all zeros
  if result == 0.0:
    return weights
  # return normalized vector (unit norm)
  return weights / result
```

```
# grid search weights
def grid search(members, testX, testy):
 # define weights to consider
 W = [0.0, 0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
 best score, best weights = 0.0, None
 # iterate all possible combinations (cartesian product)
 for weights in product(w, repeat=len(members)):
   # skip if all weights are equal
   if len(set(weights)) == 1:
      continue
   # hack, normalize weight vector
   weights = normalize(weights)
   # evaluate weights
   score = evaluate ensemble(members, weights, testX, testy)
   if score > best score:
     best_score, best_weights = score, weights
      print('>%s %.3f' % (best weights, best score))
  return list(best weights)
# fit all models
n members = 3
members = [fit_model(trainX, trainy) for _ in range(n_members)]
# evaluate each single model on the test set
testy_enc = keras.utils.to_categorical(testy)
for i in range(n members):
 , test acc = members[i].evaluate(testX, testy enc, verbose=0)
 print('Model %d: %.3f' % (i+1, test_acc))
# evaluate averaging ensemble (equal weights)
weights = [1.0/n_members for _ in range(n_members)]
score = evaluate_ensemble(members, weights, testX, testy)
print('Equal Weights Score: %.3f' % score)
# grid search weights
weights = grid search(members, testX, testy)
score = evaluate_ensemble(members, weights, testX, testy)
print('Grid Search Weights: %s, Score: %.3f' % (weights, score))
     Model 1: 0.833
     Model 2: 0.838
     Model 3: 0.832
     Equal Weights Score: 0.837
     >[0. 0. 1.] 0.832
     >[0. 1. 0.] 0.838
     >[0. 0.75 0.25] 0.840
                  0.88888889 0.11111111] 0.841
     Grid Search Weights: [0.0, 0.88888888888889, 0.1111111111111111], Score: 0.841
```

Weighted Average MLP Ensemble

```
# global optimization to find coefficients for weighted ensemble on blobs problem
from sklearn.datasets import make blobs
from sklearn.metrics import accuracy score
from keras.utils import to categorical
from keras.models import Sequential
from keras.layers import Dense
from matplotlib import pyplot
from numpy import mean
from numpy import std
from numpy import array
from numpy import argmax
from numpy import tensordot
from numpy.linalg import norm
from scipy.optimize import differential evolution
# fit model on dataset
def fit model(trainX, trainy):
  trainy enc = to categorical(trainy)
  # define model
  model = Sequential()
  model.add(Dense(25, input dim=2, activation='relu'))
  model.add(Dense(3, activation='softmax'))
 model.compile(loss='categorical_crossentropy', optimizer='adam', metrics=['accuracy'])
  # fit model
  model.fit(trainX, trainy enc, epochs=500, verbose=0)
  return model
# make an ensemble prediction for multi-class classification
def ensemble predictions(members, weights, testX):
  # make predictions
 yhats = [model.predict(testX) for model in members]
 yhats = array(yhats)
  # weighted sum across ensemble members
  summed = tensordot(yhats, weights, axes=((0),(0)))
  # argmax across classes
  result = argmax(summed, axis=1)
  return result
# evaluate a specific number of members in an ensemble
def evaluate_ensemble(members, weights, testX, testy):
  # make prediction
  yhat = ensemble predictions(members, weights, testX)
  # calculate accuracy
  return accuracy score(testy, yhat)
```

```
# normalize a vector to have unit norm
def normalize(weights):
 # calculate l1 vector norm
  result = norm(weights, 1)
  # check for a vector of all zeros
  if result == 0.0:
    return weights
  # return normalized vector (unit norm)
  return weights / result
# loss function for optimization process, designed to be minimized
def loss_function(weights, members, testX, testy):
  # normalize weights
  normalized = normalize(weights)
  # calculate error rate
  return 1.0 - evaluate ensemble(members, normalized, testX, testy)
# generate 2d classification dataset
X, y = make blobs(n samples=1100, centers=3, n features=2, cluster std=2, random state=2)
# split into train and test
n train = 100
trainX, testX = X[:n_train, :], X[n_train:, :]
trainy, testy = y[:n train], y[n train:]
print(trainX.shape, testX.shape)
# fit all models
n members = 3
members = [fit_model(trainX, trainy) for _ in range(n_members)]
# evaluate each single model on the test set
testy enc = to categorical(testy)
for i in range(n members):
  _, test_acc = members[i].evaluate(testX, testy_enc, verbose=0)
  print('Model %d: %.3f' % (i+1, test acc))
# evaluate averaging ensemble (equal weights)
weights = [1.0/n_members for _ in range(n_members)]
score = evaluate_ensemble(members, weights, testX, testy)
print('Equal Weights Score: %.3f' % score)
# define bounds on each weight
bound_w = [(0.0, 1.0) for _ in range(n_members)]
# arguments to the loss function
search arg = (members, testX, testy)
# global optimization of ensemble weights
result = differential evolution(loss function, bound w, search arg, maxiter=1000, tol=1e-7)
# get the chosen weights
weights = normalize(result['x'])
print('Optimized Weights: %s' % weights)
# evaluate chosen weights
score = evaluate ensemble(members, weights, testX, testy)
print('Optimized Weights Score: %.3f' % score)
```

(100, 2) (1000, 2) Model 1: 0.808 Model 2: 0.810 Model 3: 0.811

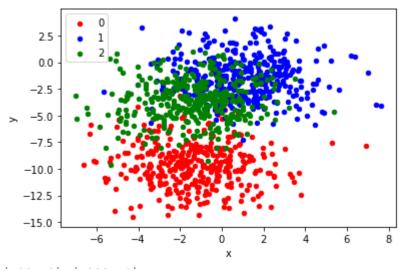
Equal Weights Score: 0.811

Optimized Weights: [0.0652782 0.35024905 0.58447275]

Optimized Weights Score: 0.816

Stacked Ensemble

```
import tensorflow as tf
from tensorflow import keras
import numpy as np
import sklearn
from sklearn.datasets import make blobs
from sklearn.linear model import LogisticRegression
from keras.utils import to categorical
from keras.models import Sequential
from keras.layers import Dense
from matplotlib import pyplot
from os import makedirs
from keras.layers.merge import concatenate
from numpy import argmax
from pandas import DataFrame
# fit model on dataset
def fit model(trainX, trainy):
  # define model
  model = Sequential()
  model.add(Dense(25, input_dim=2, activation='relu'))
  model.add(Dense(3, activation='softmax'))
  model.compile(loss='categorical_crossentropy', optimizer='adam', metrics=['accuracy'])
  # fit model
  model.fit(trainX, trainy, epochs=500, verbose=0)
  return model
# generate 2d classification dataset
X, y = make_blobs(n_samples=1100, centers=3, n_features=2, cluster_std=2, random_state=2)
df = DataFrame(dict(x=X[:,0], y=X[:,1], label=y))
colors = {0:'red', 1:'blue', 2:'green'}
fig, ax = pyplot.subplots()
grouped = df.groupby('label')
for key, group in grouped:
    group.plot(ax=ax, kind='scatter', x='x', y='y', label=key, color=colors[key])
pyplot.show()
# one hot encode output variable
y1 = to categorical(y)
# split into train and test
n train = 100
trainX, testX = X[:n train, :], X[n train:, :]
trainy, testy = y1[:n_train], y1[n_train:]
print(trainX.shape, testX.shape)
```



```
# create directory for models
#makedirs('models')
# fit and save models
n members = 5
for i in range(n_members):
  # fit model
  model = fit_model(trainX, trainy)
  # save model
  filename = 'models/model_' + str(i + 1) + '.h5'
  model.save(filename)
  print('>Saved %s' % filename)
     >Saved models/model_1.h5
     >Saved models/model 2.h5
     >Saved models/model 3.h5
     >Saved models/model 4.h5
     >Saved models/model 5.h5
# load models from file
def load_all_models(n_models):
  all models = list()
  for i in range(n_models):
    # define filename for this ensemble
    filename = 'models/model_' + str(i + 1) + '.h5'
    # load model from file
    model = keras.models.load model(filename)
    # add to list of members
    all models.append(model)
    print('>loaded %s' % filename)
  return all_models
# create stacked model input dataset as outputs from the ensemble
def stacked dataset(members, inputX):
  stackX = None
  for model in members:
    # make prediction
```

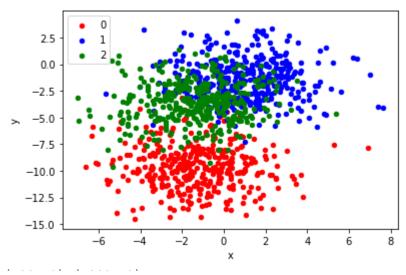
```
yhat = model.predict(inputX, verbose=0)
    # stack predictions into [rows, members, probabilities]
    if stackX is None:
      stackX = yhat
    else:
      stackX = np.dstack((stackX, yhat))
  # flatten predictions to [rows, members x probabilities]
  stackX = stackX.reshape((stackX.shape[0], stackX.shape[1]*stackX.shape[2]))
  return stackX
# fit a model based on the outputs from the ensemble members
def fit stacked model(members, inputX, inputy):
  # create dataset using ensemble
  stackedX = stacked dataset(members, inputX)
  # fit standalone model
  model = LogisticRegression()
  model.fit(stackedX, inputy)
  return model
# make a prediction with the stacked model
def stacked prediction(members, model, inputX):
  # create dataset using ensemble
  stackedX = stacked dataset(members, inputX)
  # make a prediction
 yhat = model.predict(stackedX)
  return yhat
trainy, testy = y[:n train], y[n train:]
# load all models
n members = 5
members = load all models(n members)
print('Loaded %d models' % len(members))
# evaluate standalone models on test dataset
for model in members:
  testy enc = to categorical(testy)
  _, acc = model.evaluate(testX, testy_enc, verbose=0)
  print('Model Accuracy: %.3f' % acc)
# fit stacked model using the ensemble
model = fit stacked model(members, testX, testy)
# evaluate model on test set
yhat = stacked prediction(members, model, testX)
acc = sklearn.metrics.accuracy score(testy, yhat)
print('Stacked Test Accuracy: %.3f' % acc)
     >loaded models/model 1.h5
     >loaded models/model 2.h5
     >loaded models/model 3.h5
     >loaded models/model 4.h5
     >loaded models/model 5.h5
     Loaded 5 models
```

Model Accuracy: 0.814 Model Accuracy: 0.806 Model Accuracy: 0.808 Model Accuracy: 0.809 Model Accuracy: 0.806

Stacked Test Accuracy: 0.833

Integrated Stacked Ensemble

```
import tensorflow as tf
from tensorflow import keras
import numpy as np
import sklearn
from sklearn.datasets import make blobs
from sklearn.linear model import LogisticRegression
from keras.utils import to categorical
from keras.models import Sequential
from keras.layers import Dense
from matplotlib import pyplot
from os import makedirs
from keras.layers.merge import concatenate
from numpy import argmax
from pandas import DataFrame
# fit model on dataset
def fit model(trainX, trainy):
  # define model
  model = Sequential()
  model.add(Dense(25, input_dim=2, activation='relu'))
  model.add(Dense(3, activation='softmax'))
  model.compile(loss='categorical_crossentropy', optimizer='adam', metrics=['accuracy'])
  # fit model
  model.fit(trainX, trainy, epochs=500, verbose=0)
  return model
# generate 2d classification dataset
X, y = make_blobs(n_samples=1100, centers=3, n_features=2, cluster_std=2, random_state=2)
df = DataFrame(dict(x=X[:,0], y=X[:,1], label=y))
colors = {0:'red', 1:'blue', 2:'green'}
fig, ax = pyplot.subplots()
grouped = df.groupby('label')
for key, group in grouped:
    group.plot(ax=ax, kind='scatter', x='x', y='y', label=key, color=colors[key])
pyplot.show()
# one hot encode output variable
y1 = to categorical(y)
# split into train and test
n train = 100
trainX, testX = X[:n train, :], X[n train:, :]
trainy, testy = y1[:n_train], y1[n_train:]
print(trainX.shape, testX.shape)
```



```
# create directory for models
#makedirs('models')
# fit and save models
n members = 5
for i in range(n_members):
  # fit model
  model = fit_model(trainX, trainy)
  # save model
  filename = 'models/model_' + str(i + 1) + '.h5'
  model.save(filename)
  print('>Saved %s' % filename)
     >Saved models/model_1.h5
     >Saved models/model 2.h5
     >Saved models/model 3.h5
     >Saved models/model 4.h5
     >Saved models/model 5.h5
# load models from file
def load_all_models(n_models):
  all models = list()
  for i in range(n_models):
    # define filename for this ensemble
    filename = 'models/model_' + str(i + 1) + '.h5'
    # load model from file
    model = keras.models.load model(filename)
    # add to list of members
    all models.append(model)
    print('>loaded %s' % filename)
  return all_models
# define stacked model from multiple member input models
def define stacked model(members):
  # update all layers in all models to not be trainable
  for i in range(len(members)):
    model = members[i]
```

```
for layer in model.layers:
      # make not trainable
      layer.trainable = False
      # rename to avoid 'unique layer name' issue
      layer._name = 'ensemble_' + str(i+1) + ' ' + layer.name
  # define multi-headed input
  ensemble visible = [model.input for model in members]
  # concatenate merge output from each model
  ensemble outputs = [model.output for model in members]
  merge = concatenate(ensemble outputs)
  hidden = Dense(10, activation='relu')(merge)
  output = Dense(3, activation='softmax')(hidden)
  model = keras.Model(inputs=ensemble visible, outputs=output)
  # plot graph of ensemble
  keras.utils.plot model(model, show shapes=True, to file='model graph.png')
  # compile
  model.compile(loss='categorical crossentropy', optimizer='adam', metrics=['accuracy'])
  return model
# fit a stacked model
def fit stacked model(model, inputX, inputy):
 # prepare input data
 X = [inputX for in range(len(model.input))]
  # encode output data
  inputy_enc = to_categorical(inputy)
  # fit model
 model.fit(X, inputy_enc, epochs=300, verbose=0)
# make a prediction with a stacked model
def predict stacked model(model, inputX):
  # prepare input data
 X = [inputX for in range(len(model.input))]
  # make prediction
  return model.predict(X, verbose=0)
trainy, testy = y[:n_train], y[n_train:]
# load all models
n members = 5
members = load all models(n members)
print('Loaded %d models' % len(members))
# define ensemble model
stacked model = define stacked model(members)
# fit stacked model on test dataset
fit stacked model(stacked model, testX, testy)
# make predictions and evaluate
yhat = predict stacked model(stacked model, testX)
yhat = argmax(yhat, axis=1)
acc = sklearn.metrics.accuracy score(testy, yhat)
print('Integrated Stacked Test Accuracy: %.3f' % acc)
```

```
>loaded models/model_1.h5
>loaded models/model_2.h5
>loaded models/model_3.h5
>loaded models/model_4.h5
>loaded models/model_5.h5
Loaded 5 models
Integrated Stacked Test Accuracy: 0.828
```