Setting Up the Neural Network

In the first stage (w pierwszym etapie), we define a simple architecture of the neural network and train it to recognize digits in the MNIST dataset.

```
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import matplotlib.animation
from keras.models import Sequential
from keras.layers import Dense, Dropout
from keras.optimizers import SGD
from keras import backend as K

from keras.datasets import mnist
from keras.utils import to_categorical
from sklearn.model_selection import train_test_split

%matplotlib inline
plt.rcParams["animation.html"] = "jshtml"

nb_classes = 10
```

The dropout layers have the very specific function to drop out (wykluczyć) a random set of activations in that layer by setting them to zero in the forward pass. This allows the model to avoid overfitting but should only be used at training time, not at test time.

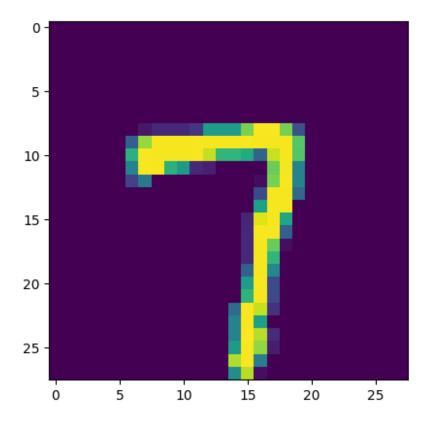
```
# Set dropout rate - fractions of neurons to drop
dropout = 0.5
# Build very simple neural network with 2 hidden layers
model = Sequential()
model.add(Dense(256, activation='relu', input shape=(784,)))
model.add(Dropout(dropout))
model.add(Dense(64, activation='relu'))
model.add(Dropout(dropout))
model.add(Dense(nb classes, activation='softmax'))
model.compile(loss='categorical crossentropy', optimizer='adam',
              metrics=['accuracy'])
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\keras\src\layers\core\dense.py:87: UserWarning: Do not pass
an `input_shape`/`input_dim` argument to a layer. When using
Sequential models, prefer using an `Input(shape)` object as the first
layer in the model instead.
  super().__init__(activity_regularizer=activity_regularizer,
**kwaras)
```

Preparing the Data

```
# The binary_crossentropy loss expects a one-hot-vector as input,
# so we apply the to categorical function from keras.utils to convert
integer labels to one-hot-vectors.
(X train, y train), (X test, y test) = mnist.load data()
X_{\text{train}} = X_{\text{train.reshape}}(60000, 784)
X \text{ test} = X \text{ test.reshape}(10000, 784)
X_{\text{train}} = \overline{X}_{\text{train.astype}}(\text{"float32"})
X test = X test.astype("float32")
# Put everything on grayscale
X train /= 255
X test /= 255
# Convert class vectors to binary class matrices
Y train = to categorical(y train, 10)
Y test = to categorical(y test, 10)
# Split training and validation data
X_train, X_val, Y_train, Y_val = train_test_split(X_train, Y_train,
train size=5/6)
```

Let's visualize an example digit:

```
# Show example digit
plt.imshow(X_train[0].reshape(28, 28))
<matplotlib.image.AxesImage at 0x23510325c90>
```



Training the Model

```
# When we have defined and compiled the model, it can be trained using
the fit function.
# We also use validation dataset to monitor validation loss and
accuracy.
network_history = model.fit(X_train, Y_train, batch_size=128,
                          epochs=20, verbose=1,
validation data=(X val, Y val))
Epoch 1/20
                    2s 3ms/step - accuracy: 0.6506 - loss:
391/391 —
1.0550 - val accuracy: 0.9346 - val_loss: 0.2259
Epoch 2/20
391/391 —
                        --- 1s 3ms/step - accuracy: 0.9034 - loss:
0.3420 - val_accuracy: 0.9536 - val_loss: 0.1658
Epoch 3/20
                        --- 1s 3ms/step - accuracy: 0.9323 - loss:
391/391 —
0.2455 - val_accuracy: 0.9597 - val_loss: 0.1385
Epoch 4/20
391/391 -
                       _____ 1s 3ms/step - accuracy: 0.9404 - loss:
0.2124 - val accuracy: 0.9613 - val loss: 0.1334
Epoch 5/20
391/391 ______ 1s 3ms/step - accuracy: 0.9468 - loss:
0.1906 - val accuracy: 0.9660 - val loss: 0.1211
Epoch 6/20
```

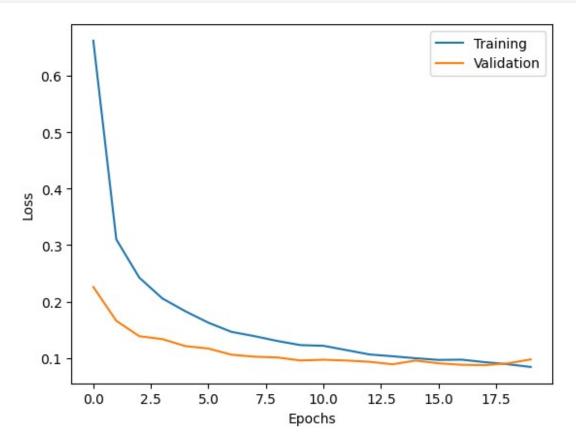
```
_____ 1s 3ms/step - accuracy: 0.9537 - loss:
391/391 ———
0.1639 - val accuracy: 0.9659 - val loss: 0.1167
Epoch 7/20
                1s 3ms/step - accuracy: 0.9584 - loss:
391/391 —
0.1465 - val accuracy: 0.9711 - val loss: 0.1058
Epoch 8/20
           _____ 1s 3ms/step - accuracy: 0.9607 - loss:
391/391 —
0.1401 - val accuracy: 0.9712 - val loss: 0.1023
0.1279 - val accuracy: 0.9723 - val loss: 0.1009
Epoch 10/20 391/391 ______ 1s 3ms/step - accuracy: 0.9651 - loss:
0.1196 - val accuracy: 0.9743 - val_loss: 0.0956
Epoch 11/20
              1s 3ms/step - accuracy: 0.9661 - loss:
391/391 ——
0.1178 - val accuracy: 0.9733 - val loss: 0.0969
Epoch 12/20
                 1s 3ms/step - accuracy: 0.9676 - loss:
391/391 —
0.1100 - val accuracy: 0.9746 - val loss: 0.0955
Epoch 13/20
                _____ 1s 3ms/step - accuracy: 0.9681 - loss:
391/391 ——
0.1060 - val accuracy: 0.9741 - val loss: 0.0933
0.0983 - val accuracy: 0.9747 - val loss: 0.0889
0.0964 - val accuracy: 0.9740 - val loss: 0.0955
0.0960 - val accuracy: 0.9766 - val loss: 0.0906
Epoch 17/20
              1s 3ms/step - accuracy: 0.9704 - loss:
391/391 ——
0.0991 - val accuracy: 0.9762 - val loss: 0.0880
Epoch 18/20
                 1s 3ms/step - accuracy: 0.9735 - loss:
391/391 ——
0.0903 - val accuracy: 0.9769 - val loss: 0.0873
Epoch 19/20
              _____ 1s 3ms/step - accuracy: 0.9745 - loss:
391/391 ——
0.0854 - val accuracy: 0.9756 - val loss: 0.0904
Epoch 20/20

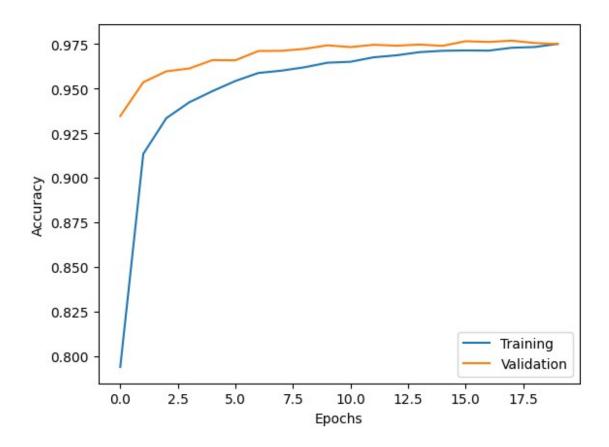
301/301 — 1s 3ms/step - accuracy: 0.9764 - loss:
0.0798 - val accuracy: 0.9750 - val loss: 0.0975
```

We can visualize the training history:

```
def plot_history(network_history):
   plt.figure()
```

```
plt.xlabel('Epochs')
    plt.ylabel('Loss')
    plt.plot(network_history.history['loss'])
    plt.plot(network history.history['val loss'])
    plt.legend(['Training', 'Validation'])
    plt.figure()
    plt.xlabel('Epochs')
    plt.ylabel('Accuracy')
    plt.plot(network history.history['accuracy'])
    plt.plot(network_history.history['val_accuracy'])
    plt.legend(['Training', 'Validation'], loc='lower right')
    plt.show()
# The fit function returns a keras.callbacks.History object which
contains the entire history
# of training/validation loss, accuracy and other metrics for each
epoch.
# We can therefore plot the behavior of loss and accuracy during the
training phase.
plot history(network history)
```





Extracting Hidden Layer Activations

```
from keras.models import Model
# Wymuś aktywację .input
_ = model.predict(X_train[:1])
# Model pomocniczy
activation model = Model(
    inputs=model.layers[0].input,
    outputs=[model.layers[0].output, model.layers[2].output,
model.layers[4].output]
# Aktywacje
layer1 output, layer2 output, layer3 output =
activation model.predict(X train)
train_ids = [np.arange(len(Y_train))[Y_train[:,i] == 1] for i in
range(10)]
1/1 -
                        0s 46ms/step
1563/1563 -
                              1s 835us/step
```

Visualizing Network Activations

The 2 visualizations below show how neuron activations work:

- 1. **First animation** shows what an example digit (number 5) looks like and what activations of neurons look like in hidden layers of the neural network.
- 2. **Second animation** checks the similarity (podobieństwo) in behavior for frames (zachowania dla klatek) showing the same digit by looking at the ensemble (zespołowe) properties. In this case, ensemble properties refers to how the neurons behave on average for a large number of frames showing the same digit.

After summing up (Po zsumowaniu) the responses of as few as 20–30 frames, the pattern in the second hidden layer is almost static. After combining (połączeniu) about 70–80 frames, the pattern in the first hidden layer also appears static. This supports the idea that only a subset of all neurons is involved in the recognition of individual digits.

This observation is important when we think about using neural networks for data visualization. We can clearly see that the activations generated by examples belonging to the same class are less chaotic than the examples themselves (one same); therefore, their visualization should give a more clustered structure.

```
from matplotlib import animation
fig, axs = plt.subplots(1, 3, figsize=(15, 3))
sample_idxs = train_ids[5][:30] # np. 30 próbek cyfrv "5"
def animate(i):
    axs[0].cla()
    axs[1].cla()
    axs[2].cla()
    img = X train[sample idxs[i]].reshape(28, 28)
    axs[0].imshow(img, cmap='gray')
    axs[0].set title("Digit Image (5)")
    axs[1].bar(range(256), layer1 output[sample idxs[i]])
    axs[1].set title("Layer 1 activations")
    axs[2].bar(range(64), layer2 output[sample idxs[i]])
    axs[2].set title("Layer 2 activations")
ani = animation.FuncAnimation(fig, animate, frames=30, interval=500)
plt.close()
from IPython.display import HTML
HTML(ani.to jshtml())
<IPython.core.display.HTML object>
```

```
fig, ax = plt.subplots(figsize=(12, 3))
sample idxs = train ids[5][:80] # 80 próbek cyfry 5
def animate avg(i):
    ax.cla()
    avg activation = layer1 output[sample idxs[:i+1]].mean(axis=0)
    ax.bar(range(256), avg_activation)
    ax.set_title(f"Average Layer 1 activation (first {i+1} samples)")
    ax.set ylim(0, 5)
ani = animation.FuncAnimation(fig, animate avg, frames=80,
interval=200)
plt.close()
HTML(ani.to jshtml())
<IPython.core.display.HTML object>
fig, ax = plt.subplots(figsize=(12, 3))
sample idxs = train ids[5][:80] # 80 próbek cyfry 5
def animate avg(i):
    ax.cla()
    avg activation = layer2 output[sample idxs[:i+1]].mean(axis=0)
    ax.bar(range(64), avg activation)
    ax.set title(f"Average Layer 2 activation (first {i+1} samples)")
    ax.set ylim(0, 10)
ani2 = animation.FuncAnimation(fig, animate avg, frames=80,
interval=200)
plt.close()
HTML(ani2.to jshtml())
<IPython.core.display.HTML object>
```

Exercise 1: MNIST Visualization Using Hidden Layer Activations

In this exercise, you will explore how the activations of hidden layers in a neural network can be used for data visualization and classification.

Task Description

- 1. Project the MNIST training data into a 2-dimensional space using different dimensionality reduction techniques:
 - t-SNE

- TriMAP
- PaCMAP
- UMAP
- 2. Use layer1_output and layer2_output to project the first and second hidden layers of the neural network into a 2-dimensional space. Apply the same dimensionality reduction methods as in step 1.
- 3. Visualize both the training and test data projections.
- 4. Use the 2-dimensional projections (2 wymiarowa projekcja) for classification:
 - Use embeddings learned (osadzeń) on raw training data (and also on hidden activations of training data) to transform test data (and hidden activations of test data) into 2-dimensional space.
 - Implement the k-nearest neighbors algorithm to classify transformed points from the test set, using points from the training set as neighbors with known class assignments.
 - Since t-SNE is a non-linear, non-parametric embedding, you can't use it to transform new points into the existing embedded space (przestrzeni osadzeń).
 Use only UMAP, which has both fit_transform (learn manifold) and transform (project new data to existing manifold (rozmaitość)) methods.
 - Try with several values of n neighbors, e.g., [3, 5, 10].
- 5. Estimate the classification accuracy using this approach:
 - Compare results for all three inputs (raw data, 1st hidden layer, 2nd hidden layer)
 - Compare results for different values of n neighbors

Code Template

Here's a template to help you get started:

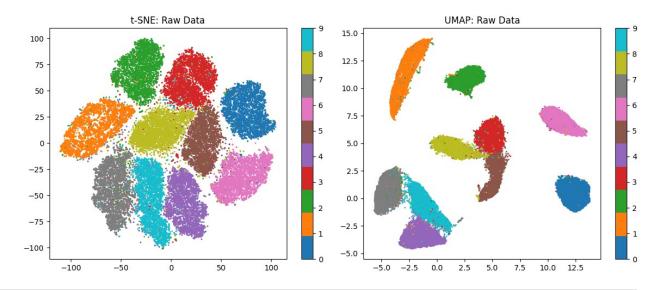
```
# To get hidden activations of test data use:
test_layer1_output, test_layer2_output, test_layer3_output =
activation_model.predict(X_test)

# To convert from one-hot encoding back to class labels:
# Use np.argmax(to_categorical(x, k), axis=1) or K.argmax
313/313 — Os 856us/step
```

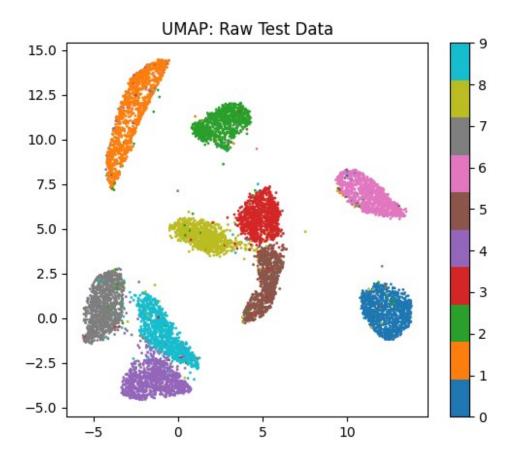
Step 1: Raw Data Embeddings Visualization

```
# Import necessary libraries
from sklearn.manifold import TSNE
import umap
# Import other visualization techniques you need
```

```
# Apply t-SNE to raw training data
tsne = TSNE(n components=2, random state=42)
X train tsne = tsne.fit transform(X train)
# Apply UMAP to raw training data
umap model = umap.UMAP(n components=\frac{2}{2}, random state=\frac{42}{2})
X train umap = umap model.fit transform(X train)
# Visualize the embeddings
plt.figure(figsize=(12, 5))
plt.subplot(1, 2, 1)
plt.scatter(X_train_tsne[:, 0], X_train_tsne[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
plt.title('t-SNE: Raw Data')
plt.colorbar()
plt.subplot(1, 2, 2)
plt.scatter(X train umap[:, 0], X train umap[:, 1],
c=np.argmax(Y_train, axis=1), cmap='tab10', s=1)
plt.title('UMAP: Raw Data')
plt.colorbar()
plt.tight layout()
plt.show()
# Apply UMAP transform to test data
X test umap = umap model.transform(X test)
# Visualize test data projection
plt.figure(figsize=(6, 5))
plt.scatter(X test umap[:, 0], X test umap[:, 1], c=np.argmax(Y test,
axis=1), cmap='tab10', s=1)
plt.title('UMAP: Raw Test Data')
plt.colorbar()
plt.show()
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\tqdm\auto.py:21: TqdmWarning: IProgress not found. Please
update jupyter and ipywidgets. See
https://ipywidgets.readthedocs.io/en/stable/user install.html
  from .autonotebook import tgdm as notebook tgdm
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force_all_finite' was renamed to 'ensure_all_finite' in 1.6 and will
be removed in 1.8.
  warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\umap\umap .py:1952: UserWarning: n jobs value 1 overridden to
1 by setting random state. Use no seed for parallelism.
  warn(
```



c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\sitepackages\sklearn\utils\deprecation.py:151: FutureWarning:
'force_all_finite' was renamed to 'ensure_all_finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(

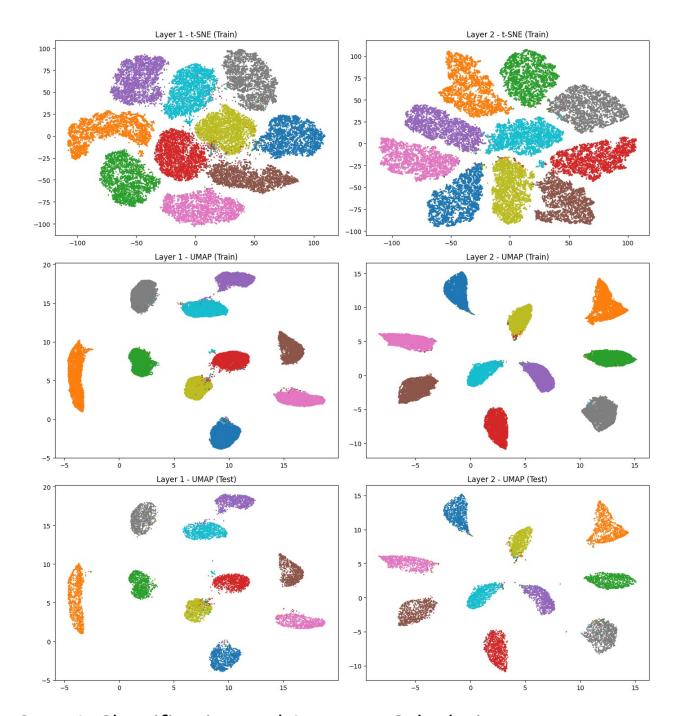


Step 2: Hidden Layer 1 Embeddings Visualization

Step 3: Hidden Layer 2 Embeddings Visualization

```
# t-SNE i UMAP dla layer1 output
tsne layer1 = TSNE(n components=2, random state=42)
X layer1 tsne = tsne layer1.fit transform(layer1 output)
umap model layer1 = umap.UMAP(n components=\frac{2}{2}, random state=\frac{42}{2})
X layer1 umap = umap model layer1.fit transform(layer1 output)
X layer1 umap test = umap model layer1.transform(test layer1 output)
# t-SNE i UMAP dla layer2 output
tsne layer2 = TSNE(n components=2, random state=42)
X layer2 tsne = tsne layer2.fit transform(layer2 output)
umap model layer2 = umap.UMAP(n components=2, random state=42)
X layer2 umap = umap model layer2.fit transform(layer2 output)
X layer2 umap test = umap model layer2.transform(test layer2 output)
# Wizualizacia
fig, axs = plt.subplots(3, 2, figsize=(14, 15))
# Layer 1 t-SNE
axs[0, 0].scatter(X layer1 tsne[:, 0], X layer1 tsne[:, 1],
c=np.argmax(Y_train, axis=1), cmap='tab10', s=1)
axs[0, 0].set title("Layer 1 - t-SNE (Train)")
# Layer 2 t-SNE
axs[0, 1].scatter(X layer2 tsne[:, 0], X layer2 tsne[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
axs[0, 1].set title("Layer 2 - t-SNE (Train)")
# Layer 1 UMAP - Train
axs[1, 0].scatter(X layer1 umap[:, 0], X layer1 umap[:, 1],
c=np.argmax(Y_train, axis=1), cmap='tab10', s=1)
axs[1, 0].set title("Layer 1 - UMAP (Train)")
# Layer 2 UMAP - Train
axs[1, 1].scatter(X_layer2_umap[:, 0], X_layer2_umap[:, 1],
c=np.argmax(Y_train, axis=1), cmap='tab10', s=1)
axs[1, 1].set title("Layer 2 - UMAP (Train)")
# Layer 1 UMAP - Test
axs[2, 0].scatter(X layer1 umap test[:, 0], X layer1 umap test[:, 1],
c=np.argmax(Y test, axis=1), cmap='tab10', s=1)
axs[2, 0].set_title("Layer 1 - UMAP (Test)")
```

```
# Layer 2 UMAP - Test
axs[2, 1].scatter(X layer2 umap test[:, 0], X layer2 umap test[:, 1],
c=np.argmax(Y test, axis=1), cmap='tab10', s=1)
axs[2, 1].set title("Layer 2 - UMAP (Test)")
plt.tight layout()
plt.show()
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force all finite' was renamed to 'ensure all finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\umap\umap .py:1952: UserWarning: n jobs value 1 overridden to
1 by setting random state. Use no seed for parallelism.
 warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force_all_finite' was renamed to 'ensure_all_finite' in 1.6 and will
be removed in 1.8.
  warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force all finite' was renamed to 'ensure all finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\umap\umap_.py:1952: UserWarning: n_jobs value 1 overridden to
1 by setting random state. Use no seed for parallelism.
  warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force all finite' was renamed to 'ensure all finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(
```



Step 4: Classification and Accuracy Calculation

```
from sklearn.neighbors import KNeighborsClassifier
from sklearn.metrics import accuracy_score

# Define a function to evaluate KNN on different embeddings
def evaluate_knn(X_train_embedded, X_test_embedded, y_train, y_test,
n_neighbors_list):
    results = []
    for n_neighbors in n_neighbors_list:
```

```
knn = KNeighborsClassifier(n neighbors=n neighbors)
        knn.fit(X train embedded, y train)
        y pred = knn.predict(X test embedded)
        accuracy = accuracy score(y test, y pred)
        results.append((n neighbors, accuracy))
        print(f"n_neighbors={n_neighbors}, Accuracy: {accuracy:.4f}")
    return results
# Convert Y train and Y test to class labels
y train = np.argmax(Y train, axis=1)
y test = np.argmax(Y test, axis=1)
# Test KNN with different numbers of neighbors
n neighbors list = [3, 5, 10]
print("Raw Data UMAP Embeddings:")
raw_results = evaluate_knn(X_train_umap, X_test_umap, y_train, y_test,
n_neighbors_list)
print("\nHidden Layer 1 UMAP Embeddings:")
layer1_results = evaluate_knn(X_layer1_umap, X layer1 umap test,
y train, y test, n neighbors list)
print("\nHidden Layer 2 UMAP Embeddings:")
layer2_results = evaluate_knn(X_layer2_umap, X_layer2_umap_test,
y train, y test, n neighbors list)
Raw Data UMAP Embeddings:
n neighbors=3, Accuracy: 0.9518
n neighbors=5, Accuracy: 0.9548
n neighbors=10, Accuracy: 0.9557
Hidden Layer 1 UMAP Embeddings:
n neighbors=3, Accuracy: 0.9697
n neighbors=5, Accuracy: 0.9714
n neighbors=10, Accuracy: 0.9713
Hidden Layer 2 UMAP Embeddings:
n neighbors=3, Accuracy: 0.9770
n neighbors=5, Accuracy: 0.9777
n neighbors=10, Accuracy: 0.9779
```

Expected Outcomes

- 1. You should observe that the visualization of hidden layer activations shows better clustering compared to raw data visualization.
- 2. The second hidden layer activations should generally show clearer class separation than the first hidden layer.

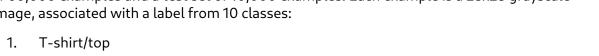
- 3. Classification accuracy should generally improve when using hidden layer activations compared to raw data, with the later layers typically providing better performance.
- 4. The choice of n_neighbors will impact performance, with the optimal value possibly differing between raw data and hidden layer embeddings.

Exercise 2: Fashion-MNIST Visualization

In this exercise, you will repeat the procedures and visualizations from Exercise 1, but now using the Fashion-MNIST dataset (or any other dataset of your choice).

Task Description

The Fashion-MNIST dataset is a dataset of Zalando's article images, consisting of a training set of 60,000 examples and a test set of 10,000 examples. Each example is a 28x28 grayscale image, associated with a label from 10 classes:



- 2. Trouser
- 3. Pullover
- 4. Dress
- 5. Coat
- 6. Sandal
- 7. Shirt
- 8. Sneaker
- 9. Bag
- 10. Ankle boot

Steps

- 1. Load and preprocess the Fashion-MNIST dataset.
- 2. Create and train a neural network similar to the one used for MNIST.
- 3. Extract the hidden layer activations.
- 4. Apply dimensionality reduction techniques (t-SNE, TriMAP, PaCMAP, UMAP) to visualize:
 - Raw data
 - First hidden layer activations

- Second hidden layer activations
- 5. Use UMAP embeddings for classification with KNN.
- 6. Compare the results with those obtained for the MNIST dataset.

Code Template

1. Load and preprocess the Fashion-MNIST dataset.

```
# Load the Fashion-MNIST dataset
from keras.datasets import fashion mnist
(X_train, y_train), (X_test, y_test) = fashion_mnist.load_data()
# Preprocess data
X_{train} = X_{train.reshape}(60000, 784)
X test = X test.reshape(10000, 784)
X train = X train.astype("float32")
X test = X test.astype("float32")
# Normalize data
X train /= 255
X \text{ test } /= 255
# Convert class vectors to binary class matrices
Y train = to categorical(y train, 10)
Y test = to categorical(y test, 10)
# Split training and validation data
X train, X val, Y train, Y val = train test split(X train, Y train,
train size=\frac{5}{6})
```

2. Create and train a neural network similar to the one used for MNIST.

```
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\keras\src\layers\core\dense.py:87: UserWarning: Do not pass
an `input shape`/`input dim` argument to a layer. When using
Sequential models, prefer using an `Input(shape)` object as the first
layer in the model instead.
 super(). init (activity regularizer=activity regularizer,
**kwargs)
391/391 ———
              ______ 2s 4ms/step - accuracy: 0.5945 - loss:
1.1321 - val accuracy: 0.8315 - val loss: 0.4755
0.5757 - val accuracy: 0.8410 - val loss: 0.4397
Epoch 3/20 ______ 1s 3ms/step - accuracy: 0.8195 - loss:
0.5152 - val accuracy: 0.8531 - val loss: 0.4005
Epoch 4/20
         _____ 1s 3ms/step - accuracy: 0.8316 - loss:
391/391 ——
0.4774 - val accuracy: 0.8590 - val_loss: 0.3865
Epoch 5/20
              _____ 1s 3ms/step - accuracy: 0.8423 - loss:
391/391 ——
0.4525 - val accuracy: 0.8712 - val loss: 0.3604
Epoch 6/20
               1s 3ms/step - accuracy: 0.8450 - loss:
391/391 —
0.4391 - val_accuracy: 0.8721 - val_loss: 0.3564
0.4278 - val accuracy: 0.8754 - val loss: 0.3400
0.4203 - val accuracy: 0.8781 - val loss: 0.3427
0.4077 - val accuracy: 0.8772 - val loss: 0.3353
0.3980 - val accuracy: 0.8840 - val loss: 0.3362
Epoch 11/20
               1s 3ms/step - accuracy: 0.8600 - loss:
391/391 ——
0.3935 - val accuracy: 0.8826 - val loss: 0.3258
Epoch 12/20
                 _____ 1s 3ms/step - accuracy: 0.8616 - loss:
391/391 ——
0.3851 - val_accuracy: 0.8835 - val_loss: 0.3301
0.3791 - val accuracy: 0.8766 - val loss: 0.3315
Epoch 14/20 ______ 1s 3ms/step - accuracy: 0.8656 - loss:
0.3764 - val accuracy: 0.8837 - val loss: 0.3237
Epoch 15/20
```

```
391/391 —
                     ——— 1s 3ms/step - accuracy: 0.8675 - loss:
0.3686 - val accuracy: 0.8837 - val loss: 0.3242
Epoch 16/20
                    _____ 1s 3ms/step - accuracy: 0.8694 - loss:
391/391 —
0.3583 - val accuracy: 0.8865 - val loss: 0.3186
Epoch 17/20
               _____ 1s 3ms/step - accuracy: 0.8713 - loss:
391/391 —
0.3605 - val accuracy: 0.8887 - val_loss: 0.3128
Epoch 18/20
              _____ 1s 3ms/step - accuracy: 0.8715 - loss:
391/391 ——
0.3580 - val accuracy: 0.8856 - val loss: 0.3197
Epoch 19/20
                 1s 3ms/step - accuracy: 0.8766 - loss:
391/391 ———
0.3448 - val accuracy: 0.8864 - val loss: 0.3171
Epoch 20/20
           _____ 1s 3ms/step - accuracy: 0.8749 - loss:
391/391 ——
0.3427 - val_accuracy: 0.8851 - val_loss: 0.3180
```

3. Extract the hidden layer activations.

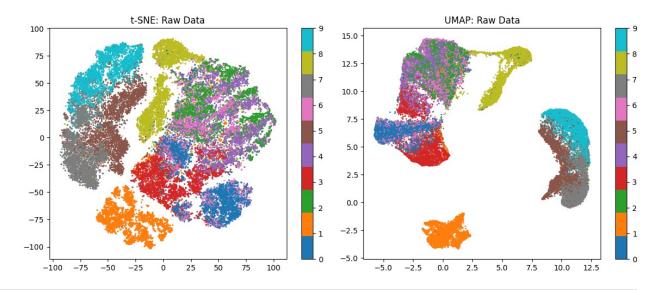
4. Apply dimensionality reduction techniques (t-SNE, TriMAP, PaCMAP, UMAP) to visualize: - Raw data - First hidden layer activations - Second hidden layer activations

```
# Apply t-SNE to raw training data
tsne = TSNE(n_components=2, random_state=42)
X_train_tsne = tsne.fit_transform(X_train)

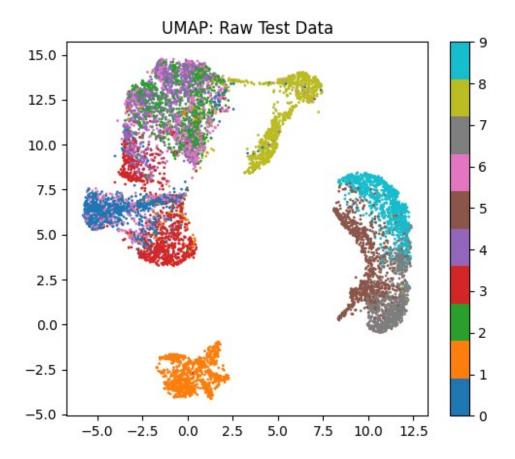
# Apply UMAP to raw training data
umap_model = umap.UMAP(n_components=2, random_state=42)
X_train_umap = umap_model.fit_transform(X_train)

# Visualize the embeddings
```

```
plt.figure(figsize=(12, 5))
plt.subplot(1, 2, 1)
plt.scatter(X_train_tsne[:, 0], X_train_tsne[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
plt.title('t-SNE: Raw Data')
plt.colorbar()
plt.subplot(1, 2, 2)
plt.scatter(X train umap[:, 0], X train umap[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
plt.title('UMAP: Raw Data')
plt.colorbar()
plt.tight layout()
plt.show()
# Apply UMAP transform to test data
X test umap = umap model.transform(X test)
# Visualize test data projection
plt.figure(figsize=(6, 5))
plt.scatter(X test umap[:, 0], X test umap[:, 1], c=np.argmax(Y test,
axis=1), cmap='tab\overline{10}', s=1)
plt.title('UMAP: Raw Test Data')
plt.colorbar()
plt.show()
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force_all_finite' was renamed to 'ensure_all_finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\umap\umap .py:1952: UserWarning: n jobs value 1 overridden to
1 by setting random state. Use no seed for parallelism.
 warn(
```

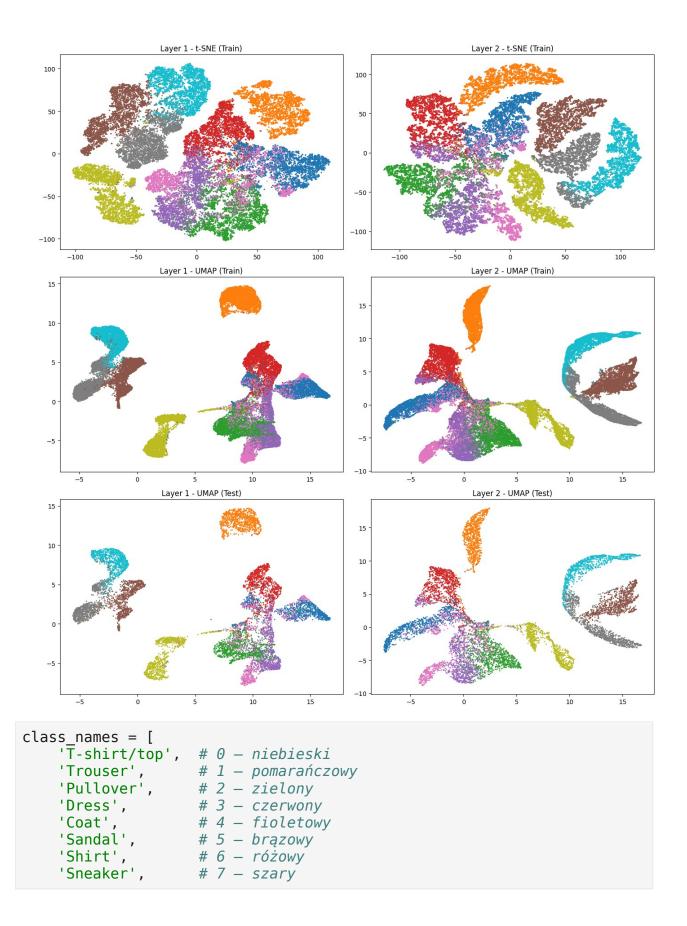


c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\sitepackages\sklearn\utils\deprecation.py:151: FutureWarning:
'force_all_finite' was renamed to 'ensure_all_finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(



```
# t-SNE i UMAP dla layer1 output
tsne layer1 = TSNE(n components=2, random state=42)
X layer1 tsne = tsne layer1.fit transform(layer1 output)
umap model layer1 = umap.UMAP(n components=2, random state=42)
X layer1 umap = umap model layer1.fit transform(layer1 output)
X layer1 umap test = umap model layer1.transform(test layer1 output)
# t-SNE i UMAP dla layer2 output
tsne layer2 = TSNE(n components=2, random state=42)
X layer2 tsne = tsne layer2.fit transform(layer2 output)
umap model layer2 = umap.UMAP(n components=\frac{2}{2}, random state=\frac{42}{2})
X layer2 umap = umap model layer2.fit transform(layer2 output)
X layer2 umap test = umap model layer2.transform(test layer2 output)
# Wizualizacia
fig, axs = plt.subplots(3, 2, figsize=(14, 15))
# Layer 1 t-SNE
axs[0, 0].scatter(X_layer1_tsne[:, 0], X_layer1_tsne[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
axs[0, 0].set title("Layer 1 - t-SNE (Train)")
# Layer 2 t-SNE
axs[0, 1].scatter(X layer2 tsne[:, 0], X layer2 tsne[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
axs[0, 1].set title("Layer 2 - t-SNE (Train)")
# Laver 1 UMAP - Train
axs[1, 0].scatter(X layer1 umap[:, 0], X layer1 umap[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
axs[1, 0].set title("Layer 1 - UMAP (Train)")
# Layer 2 UMAP - Train
axs[1, 1].scatter(X layer2 umap[:, 0], X layer2 umap[:, 1],
c=np.argmax(Y train, axis=1), cmap='tab10', s=1)
axs[1, 1].set title("Layer 2 - UMAP (Train)")
# Layer 1 UMAP - Test
axs[2, 0].scatter(X_layer1_umap_test[:, 0], X_layer1_umap_test[:, 1],
c=np.argmax(Y_test, axis=1), cmap='tab10', s=1)
axs[2, 0].set title("Layer 1 - UMAP (Test)")
# Layer 2 UMAP - Test
axs[2, 1].scatter(X layer2 umap test[:, 0], X layer2 umap test[:, 1],
c=np.argmax(Y_test, axis=1), cmap='tab10', s=1)
axs[2, 1].set_title("Layer 2 - UMAP (Test)")
```

```
plt.tight layout()
plt.show()
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force all finite' was renamed to 'ensure all finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\umap\umap .py:1952: UserWarning: n jobs value 1 overridden to
1 by setting random state. Use no seed for parallelism.
 warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force all finite' was renamed to 'ensure all finite' in 1.6 and will
be removed in 1.8.
  warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force all finite' was renamed to 'ensure all finite' in 1.6 and will
be removed in 1.8.
  warnings.warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\umap\umap .py:1952: UserWarning: n jobs value 1 overridden to
1 by setting random state. Use no seed for parallelism.
  warn(
c:\Users\jacek\AppData\Local\Programs\Python\Python311\Lib\site-
packages\sklearn\utils\deprecation.py:151: FutureWarning:
'force_all_finite' was renamed to 'ensure_all_finite' in 1.6 and will
be removed in 1.8.
 warnings.warn(
```



```
'Bag', # 8 — oliwkowy
'Ankle boot' # 9 — jasnoniebieski
]
```

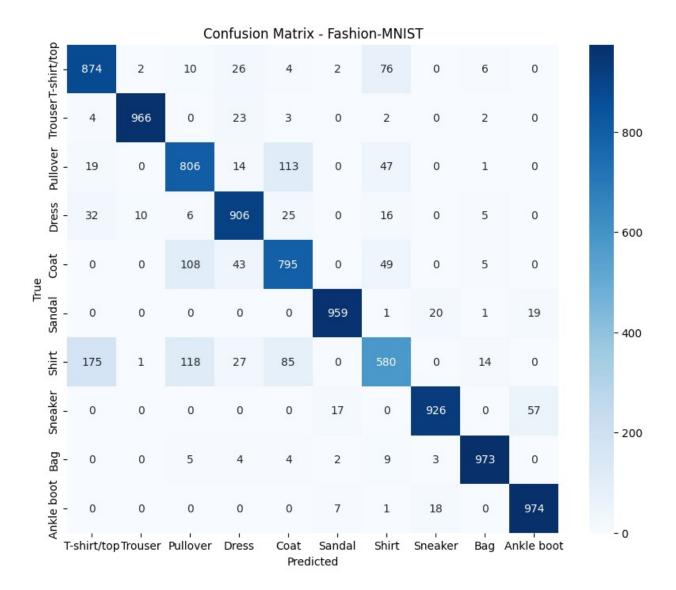
5. Use UMAP embeddings for classification with KNN.

```
# Define a function to evaluate KNN on different embeddings
def evaluate knn(X train embedded, X test embedded, y train, y test,
n_neighbors_list):
    results = []
    for n neighbors in n neighbors list:
        knn = KNeighborsClassifier(n neighbors=n neighbors)
        knn.fit(X train embedded, y train)
        y_pred = knn.predict(X_test_embedded)
        accuracy = accuracy score(y test, y pred)
        results.append((n_neighbors, accuracy))
        print(f"n neighbors={n_neighbors}, Accuracy: {accuracy:.4f}")
    return results
# Convert Y_train and Y_test to class labels
y_train = np.argmax(Y_train, axis=1)
y test = np.argmax(Y test, axis=1)
# Test KNN with different numbers of neighbors
n neighbors list = [3, 5, 10]
print("Raw Data UMAP Embeddings:")
raw results = evaluate knn(X train umap, X test umap, y train, y test,
n neighbors list)
print("\nHidden Layer 1 UMAP Embeddings:")
layer1_results = evaluate_knn(X_layer1_umap, X_layer1_umap_test,
y train, y test, n neighbors list)
print("\nHidden Layer 2 UMAP Embeddings:")
layer2 results = evaluate knn(X layer2 umap, X layer2 umap test,
y train, y test, n neighbors list)
Raw Data UMAP Embeddings:
n neighbors=3, Accuracy: 0.7415
n neighbors=5, Accuracy: 0.7596
n neighbors=10, Accuracy: 0.7721
Hidden Layer 1 UMAP Embeddings:
n neighbors=3, Accuracy: 0.8427
n neighbors=5, Accuracy: 0.8506
n_neighbors=10, Accuracy: 0.8573
```

6. Compare the results with those obtained for the MNIST dataset.

We can see that when we embed the raw data using UMAP, the classification results are worse compared to using the preprocessed data from the first and second hidden layers of the network on the Fashion-MNIST dataset. In general, we obtain better classification results from the second hidden layer. Additionally, when we test KNN with k=3,5,10, the classification accuracy increases with the value of k. The same conclusions were observed for the MNIST dataset. Furthermore, the UMAP visualizations show that the data is more separable after passing through the network layers compared to the raw input data.

```
from sklearn.metrics import confusion matrix
import seaborn as sns
import matplotlib.pyplot as plt
y true = np.argmax(Y test, axis=1)
y pred probs = model fmnist.predict(X test)
y pred = np.argmax(y pred probs, axis=1)
conf matrix = confusion matrix(y true, y pred)
class_names = ['T-shirt/top', 'Trouser', 'Pullover', 'Dress', 'Coat',
               'Sandal', 'Shirt', 'Sneaker', 'Bag', 'Ankle boot']
plt.figure(figsize=(10, 8))
sns.heatmap(conf_matrix, annot=True, fmt='d', cmap='Blues',
            xticklabels=class names, yticklabels=class names)
plt.xlabel('Predicted')
plt.vlabel('True')
plt.title('Confusion Matrix - Fashion-MNIST')
plt.show()
313/313 –
                          — 0s 1ms/step
```



Expected Outcomes

1. Compare the visualization results between MNIST and Fashion-MNIST:

- Are class boundaries (granice) clearer in one dataset versus the other?
- Do hidden layer activations provide better separation in Fashion-MNIST compared to raw data?
- Which classes in Fashion-MNIST are most easily (najbardziej) confused (mylone) in the visualizations?

2. Compare classification performance:

- How does accuracy on Fashion-MNIST compare to MNIST?
- Does using hidden layer activations provide a similar boost in performance for Fashion-MNIST as it did for MNIST?
- Which classes benefit most from using hidden layer activations for classification?

1.1 Yes, the MNIST boundaries is clealer than Fashion-MNIST boundaries, because MNIST Dataset has a simpler structure.

- 1.2 Yes, hidden layer activations provide better separation in Fashion-MNIST compared to raw input data, as visible in UMAP/t-SNE visualizations.
- 1.3 The most easily confused classes are Shirt with T-shirt.
- 2.1 The accuracy on Fashion-MNIST is lower than MNIST.
- 2.2 No, in Fashion-MNIST the improvement from using hidden layer activations is smaller compared to MNIST.
- 2.3 The classes that benefit the most from hidden layer activations for classification are Sandal, Sneaker, and Ankle boot, which are clearly separated in the deeper layers.

Discussion Questions

- 1. How does the neural network's representation of fashion items differ from its representation of digits?
- 2. What might explain any differences in visualization clarity or classification performance between the two datasets?
- 3. Which dimensionality reduction technique works best for Fashion-MNIST, and is this the same as what worked best for MNIST?
- 4. How might you modify the neural network architecture to improve visualization or classification for Fashion-MNIST specifically?
- 1. The representation of Fashion-MNIST data is more complex than that of MNIST. Digits have very distinctive shapes, which is why even the raw data performs well as a classifier.
- 2. The key difference lies in data complexity: digits have simpler visual structures, while clothing items can vary significantly in details such as shape, texture, and style.
- 3. For Fashion-MNIST, UMAP provides better class separation and more compact clusters than t-SNE, especially in the deeper hidden layers.
- 4. For example: Add more hidden layers or replace dense layers with convolutional layers.