Lab-Assignment - 5

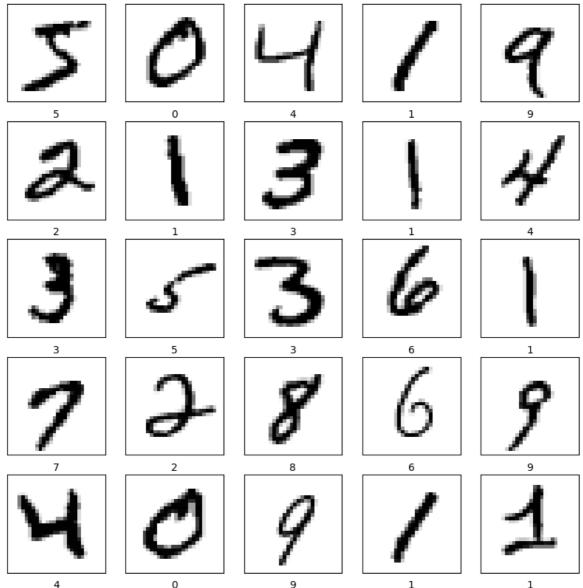
1. Load the data

```
In [1]:
          # import the libraries
          import numpy as np
          import matplotlib.pyplot as plt
          import pandas as pd
 In [2]: # Load the data
          df = pd.read_csv(r"C:\Users\raval\Downloads\mnistt\mnist_train.csv")
          df1 = pd.read_csv(r"C:\Users\raval\Downloads\mnistt\mnist_test.csv")
 Out[2]:
                 label 1x1 1x2 1x3 1x4 1x5 1x6 1x7 1x8 1x9 ... 28x19 28x20 28x21 28x22 28x23 28x24 28x2
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          60000 rows × 785 columns
In [17]: | df.info()
          <class 'pandas.core.frame.DataFrame'>
          RangeIndex: 60000 entries, 0 to 59999
          Columns: 785 entries, label to 28x28
          dtypes: int64(785)
```

2. Visualize the data columns

memory usage: 359.3 MB

```
In [14]: import numpy as np
         import matplotlib.pyplot as plt
         from tensorflow.keras.datasets import mnist # You can use TensorFlow/Keras to Load MNIST
         # Load the MNIST dataset
         (train_images, train_labels), (test_images, test_labels) = mnist.load_data()
         # Visualize the data
         plt.figure(figsize=(10, 10)) # Set the figure size for the plots
         # Plot a few sample images
         for i in range(25): # Plot the first 25 images
             plt.subplot(5, 5, i + 1) # Create a 5x5 grid of subplots
             plt.xticks([]) # Remove x-axis ticks
             plt.yticks([]) # Remove y-axis ticks
             plt.grid(False) # Turn off grid lines
             plt.imshow(train_images[i], cmap=plt.cm.binary) # Display the image
             plt.xlabel(train_labels[i]) # Add the Label as the xLabel
         plt.show() # Show the plot
```



```
In [10]: df.label.value_counts()
Out[10]: 1
              6742
              6265
         3
              6131
         2
              5958
         9
              5949
         0
              5923
         6
              5918
         8
              5851
         4
              5842
         5
              5421
         Name: label, dtype: int64
In [11]: value_counts = df['label'].value_counts()
         # Calculate the total count
         total\_count = len(df)
         # Calculate and print the percentages
         percentages = (value_counts / total_count) * 100
         # Display the result
         result_df = pd.DataFrame({'Value Counts': value_counts, 'Percentages (%)': percentages})
         result_df
Out[11]:
```

	Value Counts	Percentages (%)
1	6742	11.236667
7	6265	10.441667
3	6131	10.218333
2	5958	9.930000
9	5949	9.915000
0	5923	9.871667
6	5918	9.863333
8	5851	9.751667
4	5842	9.736667
5	5421	9.035000

no its not imbalced dataset

4. Compare leave-one-out with 2,3,5 and 10 -folds cross validation

```
In [23]: import numpy as np
         from sklearn.datasets import fetch openml
         from sklearn.model selection import cross val score
         from sklearn.linear_model import LogisticRegression
         # Load the MNIST dataset
         mnist = fetch_openml("mnist_784")
         X, y = mnist["data"], mnist["target"]
         # Convert labels to integers
         y = y.astype(np.uint8)
         # Split the dataset into training and testing sets if needed
         # You can use train_test_split from sklearn
         # Define the Logistic regression model (you can use a different model)
         model = LogisticRegression(max iter=100)
         # List of fold values to compare
         fold values = [2, 3, 5, 10, len(X)] # 2-Fold, 3-Fold, 5-Fold, 10-Fold, and Leave-One-Out
         for folds in fold values:
             # Perform cross-validation and get accuracy scores
             scores = cross val score(model, X, y, cv=folds, scoring="accuracy")
             if folds == len(X):
                  fold_name = "Leave-One-Out"
             else:
                  fold name = f"{folds}-Fold"
             print(f"{fold name} Cross Validation Accuracy: {np.mean(scores):.4f}")
         C:\python311\Lib\site-packages\sklearn\datasets\_openml.py:932: FutureWarning: The def
         ault value of `parser` will change from `'liac-arff'` to `'auto'` in 1.4. You can set `parser='auto'` to silence this warning. Therefore, an `ImportError` will be raised fr
         om 1.4 if the dataset is dense and pandas is not installed. Note that the pandas parse
         r may return different data types. See the Notes Section in fetch_openml's API doc for
         details.
           warn(
         C:\python311\Lib\site-packages\sklearn\linear_model\_logistic.py:458: ConvergenceWarni
         ng: lbfgs failed to converge (status=1):
         STOP: TOTAL NO. of ITERATIONS REACHED LIMIT.
         Increase the number of iterations (max_iter) or scale the data as shown in:
             https://scikit-learn.org/stable/modules/preprocessing.html (https://scikit-learn.o
         rg/stable/modules/preprocessing.html)
         Please also refer to the documentation for alternative solver options:
              https://scikit-learn.org/stable/modules/linear model.html#logistic-regression (htt
         ps://scikit-learn.org/stable/modules/linear_model.html#logistic-regression)
           n_iter_i = _check_optimize_result(
         C:\python311\Lib\site-packages\sklearn\linear_model\_logistic.py:458: ConvergenceWarni
```

Leave-One-Out cross-validation (LOOCV) is not suitable for datasets with a large number of samples, such as the MNIST dataset with 70,000 samples. LOOCV would create as many folds as there are samples (in this case, 70,000 folds), which is impractical and computationally expensive.

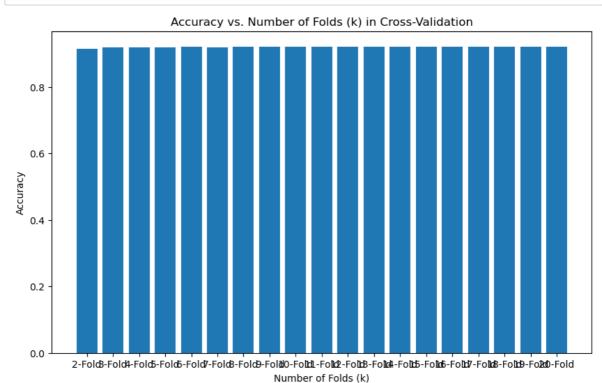
For large datasets like MNIST, it's more common to use k-fold cross-validation with a smaller value of k (e.g., 5 or 10).

- 1. 2-Fold Cross Validation Accuracy: 0.9159
- 2. 3-Fold Cross Validation Accuracy: 0.9186
- 3. 5-Fold Cross Validation Accuracy: 0.9201
- 4. 10-Fold Cross Validation Accuracy: 0.9210

These results suggest that the logistic regression model performs consistently well on the MNIST dataset, and the performance slightly improves as you increase the number of folds. It's important to note that the choice of the number of folds can affect the estimated performance, and typically, a higher number of folds provides a more robust estimate but may require more computation.

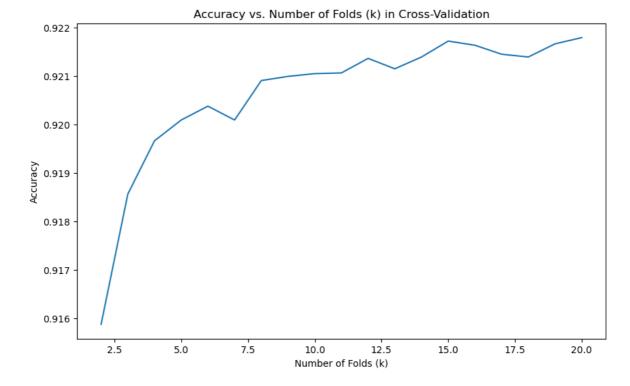
- 5. Observe the pattern as the number of folds increases
- 6. identify the good value of K for better accurracy

```
In [20]: import warnings
         warnings.filterwarnings("ignore")
         import numpy as np
         import matplotlib.pyplot as plt
         from sklearn.datasets import fetch_openml
         from sklearn.model_selection import cross_val_score
         from sklearn.linear_model import LogisticRegression
         # Load the MNIST dataset
         mnist = fetch openml("mnist 784")
         X, y = mnist["data"], mnist["target"]
         # Convert labels to integers
         y = y.astype(np.uint8)
         # Define the logistic regression model (you can use a different model)
         model = LogisticRegression(max_iter=100)
         # Lists to store fold values and accuracy scores
         fold_values = list(range(2, 21))
         accuracy_scores = []
         for folds in fold values:
             # Perform cross-validation and get accuracy scores
             scores = cross_val_score(model, X, y, cv=folds, scoring="accuracy")
             accuracy_scores.append(np.mean(scores))
         # Plot the accuracy scores for each value of k
         plt.figure(figsize=(10, 6))
         plt.bar(fold\_values, accuracy\_scores, tick\_label=[f"\{k\}-Fold" \ for \ k \ in \ fold\_values])
         plt.xlabel("Number of Folds (k)")
         plt.ylabel("Accuracy")
         plt.title("Accuracy vs. Number of Folds (k) in Cross-Validation")
         plt.show()
```



```
accuracy_scores
In [21]:
Out[21]: [0.9158714285714286,
           0.9185571609869072,
          0.919657142857143,
          0.9200857142857142,
          0.9203715802824922,
          0.9200857142857143,
           0.9209,
           0.9209859659146681,
           0.9210428571428573,
           0.9210572848611073,
           0.9213571790761814,
           0.9211430848762342,
           0.9213857142857141,
           0.9217145900392333,
          0.9216285714285715,
           0.9214431920270327,
           0.9213860460835746,
           0.9216572522888106,
           0.9217857142857142]
```

In [23]: # Plot the accuracy scores for each value of k plt.figure(figsize=(10, 6)) plt.plot(fold_values, accuracy_scores) # plt.plot(fold_values, accuracy_scores, tick_label=[f"{k}-Fold" for k in fold_values]) plt.xlabel("Number of Folds (k)") plt.ylabel("Accuracy") plt.title("Accuracy vs. Number of Folds (k) in Cross-Validation") plt.show()



Some observations from the pattern:

- 1. As k initially increases from 2 to 4, the accuracy generally improves, indicating that more folds (i.e., smaller validation sets) can lead to better model generalization.
- 2. After reaching a peak around k = 4 or k = 5, the accuracy values tend to stabilize and remain relatively consistent as k continues to increase. This suggests that further increasing the number of folds does not significantly impact model performance.
- 3. The accuracy values fluctuate slightly as k increases, but the overall trend is relatively flat, with accuracy values around 0.921.

	4. There isn't a substantial drop in accuracy even with a high number of folds (e.g., k = 21), indicating that the	
	In practice, it's common to use a moderate number of folds (e.g., 5 or 10) in cross-validation, as this strikes a balance between obtaining a reliable estimate of model performance and computational efficiency.	
In []:		