



KESHAV MEMORIAL INSTITUTE OF TECHNOLOGY
AN AUTONOMOUS INSTITUTION - ACCREDITED BY NAAC WITH 'A' GRADE
Narayanaguda, Hyderabad.

EMBEDDED LEARNING

DAY1 EXERCISE

ML MODEL VS DL MODEL

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MACHINE LEARNING MODEL

1. Importing Libraries

```
import pandas as pd
from sklearn.datasets import load_breast_cancer
from sklearn.model_selection import train_test_split
from sklearn.preprocessing import StandardScaler
from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import accuracy_score
import seaborn as sns
```

`pandas` (`pd`) is used for data manipulation and analysis.

`load_breast_cancer` from `sklearn.datasets` loads the Breast Cancer dataset.

`train_test_split` is used to split the dataset into training and testing sets.

`StandardScaler` is typically used for scaling data (though it's not used here).

`DecisionTreeClassifier` is used to create a decision tree model for classification.

`accuracy_score` is used to calculate the accuracy of the model.

`seaborn` (`sns`) is used for visualization, specifically for generating heatmaps (in this case, for the confusion matrix).

2. Loading the Breast Cancer Dataset

```
# Load Breast Cancer dataset  
data = load_breast_cancer()
```

This line loads the Breast Cancer dataset from `sklearn.datasets`. The dataset contains 30 features describing cell characteristics of breast cancer tumors and a target variable (`y`) indicating whether the tumor is malignant or benign.

3. Extracting Features and Target Variables

```
# Load features and target  
X = data.data  
y = data.target
```

`X` contains the feature matrix (30 numerical features describing the tumors).

`y` contains the target vector, where 1 indicates malignant tumors and 0 indicates benign tumors.

4. Converting the Dataset to a DataFrame and Printing First Few Rows

```
# Convert to DataFrame for better visualization
df = pd.DataFrame(X, columns=data.feature_names)
df['target'] = y

# Print the first few rows of the dataset
print("Dataset (First 5 Rows):")
print(df.head())
```

The feature matrix (`X`) is converted into a **pandas DataFrame** with **column names** taken from `data.feature_names`.

A new column, `'target'`, is added to the DataFrame to store the target labels (`y`), allowing for better visualization and manipulation.

Printing the first few rows of the dataset (`df.head()`) helps to quickly inspect the structure of the data, including both the features and the target variable.

Dataset (First 5 Rows):

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	\
0	17.99	10.38	122.80	1001.0	0.11840	
1	20.57	17.77	132.90	1326.0	0.08474	
2	19.69	21.25	130.00	1203.0	0.10960	
3	11.42	20.38	77.58	386.1	0.14250	
4	20.29	14.34	135.10	1297.0	0.10030	

	mean compactness	mean concavity	mean concave points	mean symmetry	\
0	0.27760	0.3001	0.14710	0.2419	
1	0.07864	0.0869	0.07017	0.1812	
2	0.15990	0.1974	0.12790	0.2069	
3	0.28390	0.2414	0.10520	0.2597	
4	0.13280	0.1980	0.10430	0.1809	

	mean fractal dimension	...	worst texture	worst perimeter	worst area	\
0	0.07871	...	17.33	184.60	2019.0	
1	0.05667	...	23.41	158.80	1956.0	
2	0.05999	...	25.53	152.50	1709.0	
3	0.09744	...	26.50	98.87	567.7	
4	0.05883	...	16.67	152.20	1575.0	

	worst smoothness	worst compactness	worst concavity	worst concave points	\
0	0.1622	0.6656	0.7119	0.2654	
1	0.1238	0.1866	0.2416	0.1860	
2	0.1444	0.4245	0.4504	0.2430	
3	0.2098	0.8663	0.6869	0.2575	
4	0.1374	0.2050	0.4000	0.1625	

	worst symmetry	worst fractal dimension	target
0	0.4601	0.11890	0
1	0.2750	0.08902	0
2	0.3613	0.08758	0
3	0.6638	0.17300	0
4	0.2364	0.07678	0

[5 rows x 31 columns]

5. Splitting the Data into Training and Testing Sets

```
# Split into training and testing sets  
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)
```

The data is split into **training (80%)** and **testing (20%)** sets using `train_test_split`.

`random_state=42` ensures the results are reproducible.

6. Initializing the Decision Tree Classifier

```
# Initialize Decision Tree Classifier  
dt_model = DecisionTreeClassifier(random_state=42)
```

A **Decision Tree Classifier** is created and initialized with a fixed `random_state=42` to ensure consistent results.

7. Training the Decision Tree Model

```
# Train the model  
dt_model.fit(X_train, y_train)
```

The Decision Tree model is trained using the training data (X_train, y_train).

8. Making Predictions

```
# Predict on the test data  
y_pred_dt = dt_model.predict(X_test)
```

The trained model is used to predict the labels for the **test data** (`X_test`).

9. Evaluating the Model

```
# Evaluate the model  
dt_accuracy = accuracy_score(y_test, y_pred_dt)
```

The **accuracy** of the model is calculated by comparing the predicted labels (`y_pred_dt`) with the actual test labels (`y_test`).

10. Printing the Accuracy

```
print("Decision Tree Results:")  
print(f"Accuracy: {dt_accuracy:.2f}")
```

The **accuracy** of the Decision Tree model is printed, formatted to two decimal places.

11. Visualizing the Confusion Matrix

```
# Visualization 1: Confusion Matrix for Decision Tree
plt.figure(figsize=(8, 6))
sns.heatmap(dt_conf_matrix, annot=True, fmt='d', cmap='Blues', xticklabels=data.target_names, yticklabels=data.target_names)
plt.title('Decision Tree Confusion Matrix')
plt.xlabel('Predicted')
plt.ylabel('Actual')
plt.show()
```

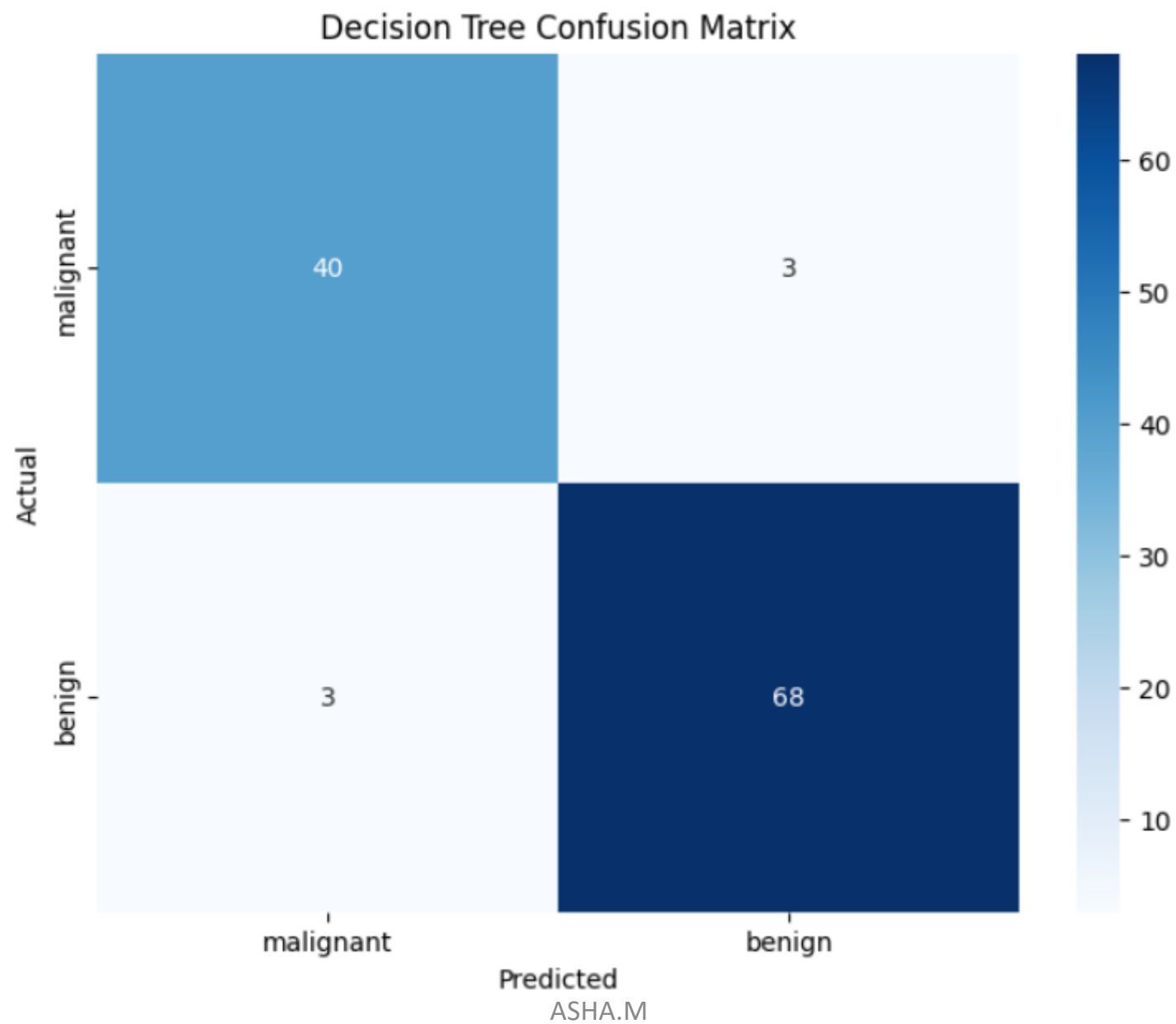
The **Confusion Matrix** is meant to be visualized using **seaborn's heatmap**.

However, the confusion matrix (`dt_conf_matrix`) needs to be computed using `confusion_matrix(y_test, y_pred_dt)`.

Once the confusion matrix is computed, `sns.heatmap` can be used to visualize it, with `xticklabels` and `yticklabels` set to the target names (`malignant` and `benign`).

Decision Tree Results:

Accuracy: 0.95



Summary of the Process:

- 1. Data Loading & Preparation:** Load the Breast Cancer dataset and convert it into a pandas DataFrame for better visualization and manipulation.
- 2. Train-Test Split:** Split the dataset into training and testing sets.
- 3. Model Initialization & Training:** Initialize and train a Decision Tree classifier.
- 4. Prediction & Evaluation:** Make predictions and evaluate the model using accuracy.
- 5. Visualization:** Convert the dataset into a DataFrame and print the first few rows to understand the structure of the data. A confusion matrix is also visualized (though it needs a fix).

DEEP LEARNING MODEL

1. Importing Libraries

```
import pandas as pd
from sklearn.datasets import load_breast_cancer
from sklearn.model_selection import train_test_split
from sklearn.preprocessing import StandardScaler
from sklearn.metrics import accuracy_score
import tensorflow as tf
from tensorflow.keras.models import Sequential
from tensorflow.keras.layers import Dense
```

`pandas` : A library used to handle data in tabular form (DataFrames).

`sklearn.datasets.load_breast_cancer` : A function that loads the Breast Cancer dataset.

`train_test_split` : Splits the data into training and test sets.

`StandardScaler` : Used to standardize the features (making them have zero mean and unit variance).

`accuracy_score` : A function that calculates the accuracy of the model.

`tensorflow` : A machine learning framework for building and training neural networks.

`Sequential` : Used to define a neural network in a linear fashion (adding layers one by one).

`Dense` : A fully connected neural network layer, used to define the number of neurons in each layer.

2. Loading the Breast Cancer Dataset

```
# Load Breast Cancer dataset  
data = load_breast_cancer()
```

This line loads the Breast Cancer dataset from `sklearn.datasets`. The dataset contains 30 features describing cell characteristics of breast cancer tumors and a target variable (`y`) indicating whether the tumor is malignant or benign.

3. Extracting Features and Target Variables

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# Load features and target  
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`X` contains the feature matrix (30 numerical features describing the tumors).

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The feature matrix (`X`) is converted into a **pandas DataFrame** with **column names** taken from `data.feature_names`.

A new column, `'target'`, is added to the DataFrame to store the target labels (`y`), allowing for better visualization and manipulation.

Printing the first few rows of the dataset (`df.head()`) helps to quickly inspect the structure of the data, including both the features and the target variable.

Dataset (First 5 Rows):

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	worst symmetry	worst fractal dimension	target
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1	0.2750	0.08902	0
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4	0.2364	0.07678	0

[5 rows x 31 columns]

5. Splitting the Data into Training and Testing Sets

```
# Split into training and testing sets  
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)
```

The data is split into **training (80%)** and **testing (20%)** sets using `train_test_split`.

`random_state=42` ensures the results are reproducible.

6. Standardizing the Features

```
# Standardize features for ANN model
scaler = StandardScaler()
X_train_scaled = scaler.fit_transform(X_train)
X_test_scaled = scaler.transform(X_test)
```

`StandardScaler()` standardizes the features to have zero mean and unit variance.

- `fit_transform(X_train)` scales the training data and learns the scaling parameters.
- `transform(X_test)` applies the same scaling transformation to the test data, ensuring no data leakage.

7. Building the ANN Model

```
# Initialize ANN model
ann_model = Sequential([
    Dense(64, activation='relu', input_dim=X_train_scaled.shape[1]), # Input layer
    Dense(32, activation='relu'),
    Dense(1, activation='sigmoid') # Output layer for binary classification
])
```

`Sequential()` : This is used to define the model layer by layer.

`Dense(64, activation='relu', input_dim=X_train_scaled.shape[1])` : The input layer with 64 neurons and ReLU activation. The `input_dim` is set to the number of features in the dataset (`X_train_scaled.shape[1]`).

`Dense(32, activation='relu')` : A hidden layer with 32 neurons and ReLU activation.

`Dense(1, activation='sigmoid')` : The output layer with 1 neuron and sigmoid activation for binary classification (0 or 1).

8. Compiling the Model

```
# Compile the model  
ann_model.compile(optimizer='adam', loss='binary_crossentropy', metrics=['accuracy'])
```

`optimizer='adam'` : Adam optimizer is used, which adapts the learning rate during training.

`loss='binary_crossentropy'` : This loss function is used for binary classification tasks. It measures the difference between the predicted and actual values.

`metrics=['accuracy']` : This metric is used to track the accuracy of the model during training and evaluation.

9. Training the Model

```
# Train the model  
ann_model.fit(X_train_scaled, y_train, epochs=10)
```

`fit(X_train_scaled, y_train, epochs=10)` : The model is trained using the training data (`X_train_scaled` and `y_train`) for 10 epochs (iterations over the entire dataset).

10. Making Predictions on the Test Set

```
# Predict on the test set  
y_pred_ann = (ann_model.predict(X_test_scaled) > 0.5).astype("int32")
```

`ann_model.predict(X_test_scaled)` : The model predicts probabilities (between 0 and 1) for the test data.

`(ann_model.predict(X_test_scaled) > 0.5)` : Converts the probabilities into binary predictions (1 for probabilities > 0.5, and 0 otherwise).

`.astype("int32")` : Converts the predicted values to integers (0 or 1).

11. Evaluating the Model

```
# Evaluate the model  
ann_accuracy = accuracy_score(y_test, y_pred_ann)
```

`accuracy_score(y_test, y_pred_ann)` : Calculates the accuracy by comparing the actual labels (`y_test`) with the predicted labels (`y_pred_ann`).

`print(f"Accuracy: {ann_accuracy:.2f}")` : Prints the accuracy of the model rounded to two decimal places.

Epoch 1/10

```
/usr/local/lib/python3.11/dist-packages/keras/src/layers/core/dense.py:87: UserWarning: Do not pass an `input_shape`,
    super().__init__(activity_regularizer=activity_regularizer, **kwargs)
```

15/15 ————— **1s** 2ms/step - accuracy: 0.6609 - loss: 0.6175

Epoch 2/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9496 - loss: 0.3162

Epoch 3/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9555 - loss: 0.1895

Epoch 4/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9542 - loss: 0.1561

Epoch 5/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9661 - loss: 0.1083

Epoch 6/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9685 - loss: 0.1037

Epoch 7/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9746 - loss: 0.0887

Epoch 8/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9888 - loss: 0.0612

Epoch 9/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9830 - loss: 0.0608

Epoch 10/10

15/15 ————— **0s** 2ms/step - accuracy: 0.9752 - loss: 0.0825

4/4 ————— **0s** 37ms/step

ANN Results:

Accuracy: 0.98

12. Visualizing the Confusion Matrix

```
# Visualization 2: Confusion Matrix for ANN
plt.figure(figsize=(8, 6))
sns.heatmap(ann_conf_matrix, annot=True, fmt='d', cmap='Blues', xticklabels=data.target_names, yticklabels=data.target_names)
plt.title('ANN Confusion Matrix')
plt.xlabel('Predicted')
plt.ylabel('Actual')
plt.show()
```

`sns.heatmap()` : It creates a heatmap of the confusion matrix, where the color intensity represents the count of predictions.

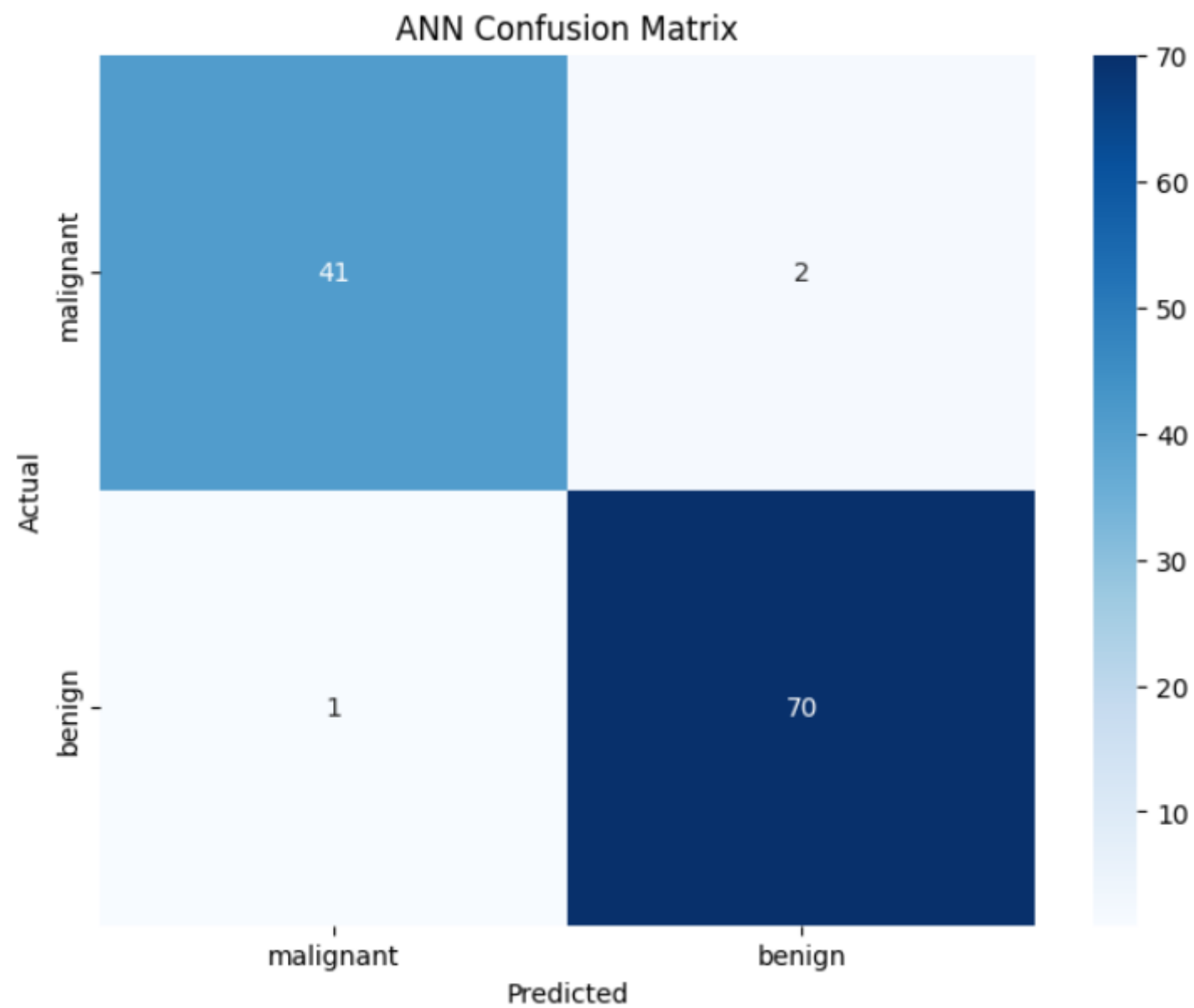
`annot=True` : Displays the actual numbers inside each matrix cell.

`fmt='d'` : Formats the numbers as integers.

`cmap='Blues'` : Uses a blue color scheme to show the matrix values.

`xticklabels` and `yticklabels` : Labels the x-axis and y-axis with the actual and predicted class names (Benign, Malignant).

`plt.show()` : Displays the heatmap.



Summary:

- 1. Data Loading:** Loads the Breast Cancer dataset and separates the features (X) and target labels (y).
- 2. Data Preprocessing:** Standardizes the features for better performance during training.
- 3. Model Building:** Constructs a neural network with 1 input layer, 1 hidden layer, and 1 output layer.
- 4. Model Training:** Trains the model for 10 epochs using the training data.
- 5. Evaluation:** Makes predictions on the test set and evaluates the accuracy of the model.

ML MODEL VS DL MODEL

1. Higher Accuracy in DL Model:

- 1.The **Deep Learning (DL) model** performs better with an accuracy of **98%**, compared to the **95%** of the **Machine Learning (ML) model**.
- 2.This indicates that the DL model is slightly more effective in making correct predictions.

2. Model Complexity:

- 1.The **ML model** (e.g., Decision Trees, Random Forest) is simpler, which means it can perform well but may not capture complex patterns in the data.
- 2.On the other hand, the **DL model** (Artificial Neural Network) has more layers and neurons, enabling it to learn and generalize better from the data, especially when the data is large and complex.

3. Data Requirements:

- 1.**ML models** generally perform well with smaller datasets and require less computational power.
- 2.**DL models**, while offering higher accuracy, typically require larger datasets and more computational resources to train effectively. However, once trained, they can outperform traditional ML models.

4. Generalization:

1. The **DL model** likely has better **generalization**, meaning it can better predict unseen data (test data) compared to the **ML model**

5. Trade-offs:

1. **ML models** are easier to train, require fewer resources, and are often faster, but may not handle very complex problems as well as **DL models**.
 2. **DL models** can achieve higher accuracy, but they are computationally expensive and may take longer to train.
- In conclusion, while the **DL model** achieves a higher accuracy, **ML models** are often more practical for simpler tasks or when computational resources are limited.