Experiment no 3:

Name:Jhanvi Parekh

Sap:60009210033

Lab exercise to be performed:

Consider fake news data set

Data Preprocessing

- 1. Load the Dataset
- 2. Text Cleaning:

Clean the text data (remove special characters, convert to lowercase, etc.).

3. Tokenization and Padding:

Tokenize and pad the text sequences.

4. Split Dataset:

Split the data into training and test sets.

- 2.Building the RNN Model
- 3. Building the LSTM Model
- 4. Building GRU Model
- 5. Perform Fake News Classification for English News

```
pata Preprocessing
import pandas as pd
import re
df = pd.read_csv('/content/fake_or_real_news 1.csv')
df.head()
```

₹		Unnamed: 0	title	text	label
	0	8476	You Can Smell Hillary's Fear	Daniel Greenfield, a Shillman Journalism Fello	FAKE
	1	10294	Watch The Exact Moment Paul Ryan Committed Pol	Google Pinterest Digg Linkedin Reddit Stumbleu	FAKE
	2	3608	Kerry to go to Paris in gesture of sympathy	U.S. Secretary of State John F. Kerry said Mon	REAL
	3	10142	Bernie supporters on Twitter erupt in anger ag	— Kaydee King (@KaydeeKing) November 9, 2016 T	FAKE
	4	875	The Battle of New York: Why This Primary Matters	It's primary day in New York and front-runners	REAL

```
#Text Cleaning
def clean_text(text):
    text = text.lower()
    text = re.sub(r'[^a-z\s]', '', text)
    text = re.sub(r'\s+', ' ', text).strip()
    return text

df['cleaned_text'] = df['text'].apply(clean_text)

df[['title', 'cleaned_text', 'label']].head()
```

	title	cleaned_text	
0	You Can Smell Hillary's Fear	daniel greenfield a shillman journalism fellow	FAKE
1	Watch The Exact Moment Paul Ryan Committed Pol	google pinterest digg linkedin reddit stumbleu	FAKE
2	Kerry to go to Paris in gesture of sympathy	us secretary of state john f kerry said monday	REAL
3	Bernie supporters on Twitter erupt in anger ag	kavdee king kavdeeking november the lesson fro	FAKE

```
CO △ ACL_Lab_3.ipynb ☆
       File Edit View Insert Runtime Tools Help <u>Last saved at 11:04 PM</u>
     + Code + Text
∷
       [ ] missing_data = df.isnull().sum()
Q
            print(missing_data)
       → Unnamed: 0
\{x\}
            title
            text
label
©∓7
            cleaned text
            dtype: int64
[ ] #Tokenization and Padding
            from tensorflow.keras.preprocessing.text import Tokenizer
            from tensorflow.keras.preprocessing.sequence import pad_sequences
       [ ] tokenizer = Tokenizer(num_words=10000)
         tokenizer.fit_on_texts(df['cleaned_text'])
       [ ] sequences = tokenizer.texts_to_sequences(df['cleaned_text'])
       [ ] #Padding
            max_sequence_length = 100
            padded sequences = pad sequences(sequences, maxlen=max sequence length)
            print(padded_sequences)
       [ 12 1 69 ... 390 745 23]
[ 9 338 1 ... 1136 1088 624]
[ 1 103 2343 ... 612 3 6648]
             [ 10 1 2311 ... 2 109 361]
[ 1 109 40 ... 3470 16 1014]
[ 311 9 784 ... 23 1 7621]]
<>
\equiv
       [ ] from sklearn.model_selection import train_test_split
>_
            y = df['label'].apply(lambda x: 1 if x == 'FAKE' else 0).values
  from sklearn.model_selection import train_test_split
      y = df['label'].apply(lambda x: 1 if x == 'FAKE' else 0).values
      X_train, X_test, y_train, y_test = train_test_split(padded_sequences, y, test_size=0.2, random_state=42)
       print(f"Training set shape: {X_train.shape}")
      print(f"Test set shape: {X_test.shape}")
      print(f"y: {y}")
  Training set shape: (5068, 100)
Test set shape: (1267, 100)
y: [1 1 0 ... 1 0 0]
 [ ] #Building the RNN Model
       from tensorflow.keras.models import Sequential
       from tensorflow.keras.layers import Embedding, SimpleRNN, Dense, Dropout
       model = Sequential()
 [ ] model.add(Embedding(input_dim=10000, output_dim=128, input_length=max_sequence_length))
      model.add(SimpleRNN(64, return_sequences=False))
       # model.add(Dropout(0.2))
      model.add(Dense(1, activation='sigmoid'))
  🚁 /usr/local/lib/python3.10/dist-packages/keras/src/layers/core/embedding.py:90: UserWarning: Argument `input_length` is deprecated. Just remove it.
         warnings.warn(
 [ ]
       model.compile(optimizer='adam',
                      loss='binary_crossentropy',
                      metrics=['accuracy'])
```



```
[ ] predictions = model.predict(X_test)
      predicted_labels = (predictions > 0.5).astype(int)
      print("Predicted Labels:")
      print(predicted_labels)
<del>→</del> 40/40 -
                                  --- 1s 12ms/step
      Predicted Labels:
      [[1]
       [1]
       ...
[0]
       [0]
       [1]]
[] for i in range(10):
         print(f"Sample {i+1}")
print(f"Sample {i+1}")
print(f"Actual Label: {'FAKE' if y_test[i] == 1 else 'REAL'}")
print(f"Predicted Label: {'FAKE' if predicted_labels[i] == 1 else 'REAL'}")
print("-" * 30)
Sample 1
Actual Label: FAKE
      Predicted Label: FAKE
      Sample 2
Actual Label: FAKE
      Predicted Label: FAKE
      Sample 3
      Actual Label: FAKE
      Predicted Label: FAKE
      Sample 4
      Actual Label: FAKE
```

```
D .....

→ Sample 6

    Actual Label: FAKE
    Predicted Label: FAKE
    Sample 7
    Actual Label: REAL
    Predicted Label: REAL
    -----
    Sample 8
    Actual Label: FAKE
    Predicted Label: FAKE
    -----
    Sample 9
    Actual Label: REAL
    Predicted Label: REAL
    -----
    Sample 10
    Actual Label: FAKE
    Predicted Label: FAKE
[ ] #Building the LSTM Model
    from tensorflow.keras.layers import Embedding, LSTM, Dense, Dropout
[ ] model1 = Sequential()
    model1.add(Embedding(input_dim=10000, output_dim=128, input_length=max_sequence_length))
    model1.add(LSTM(64, return_sequences=False))
    model1.add(Dropout(0.5))
    model1.add(Dense(1, activation='sigmoid'))
[ ] model1.compile(optimizer='adam',
                  loss='binary_crossentropy',
                  metrics=['accuracy'])
 [ ] model1.compile(optimizer='adam',
                 loss='binary_crossentropy',
                  metrics=['accuracy'])
 history1 = model1.fit(X_train, y_train,
                        epochs=5,
                       batch_size=64,
                       validation_data=(X_test, y_test))

→ Epoch 1/5

                           -- 14s 150ms/step - accuracy: 0.6191 - loss: 0.6372 - val_accuracy: 0.8579 - val_loss: 0.3389
     80/80 -
     Epoch 2/5
                           --- 12s 144ms/step - accuracy: 0.9111 - loss: 0.2452 - val_accuracy: 0.8437 - val_loss: 0.3431
     80/80 -
     Epoch 3/5
     80/80 -
                            -- 12s 147ms/step - accuracy: 0.9615 - loss: 0.1319 - val_accuracy: 0.8650 - val_loss: 0.3925
     Epoch 4/5
     80/80 -
                           --- 12s 149ms/step - accuracy: 0.9861 - loss: 0.0486 - val_accuracy: 0.8556 - val_loss: 0.4358
     Epoch 5/5
     80/80 -
                           --- 20s 143ms/step - accuracy: 0.9933 - loss: 0.0236 - val_accuracy: 0.8382 - val_loss: 0.6501
 [ ] loss, accuracy = model1.evaluate(X_test, y_test)
     print(f"Test Loss: {loss}")
     print(f"Test Accuracy: {accuracy}")
 - 1s 17ms/step - accuracy: 0.8320 - loss: 0.6778
     Test Loss: 0.6500509977340698
```

Test Accuracy: 0.8382004499435425

```
[] # Building GRU Model
     from tensorflow.keras.layers import Embedding, GRU, Dense, Dropout
[ ] model2 = Sequential()
     model2.add(Embedding(input_dim=10000, output_dim=128, input_length=max_sequence_length))
     model2.add(GRU(64, return_sequences=False))
     model2.add(Dropout(0.5))
     model2.add(Dense(1, activation='sigmoid'))
[] model2.compile(optimizer='adam',
                   loss='binary_crossentropy',
                   metrics=['accuracy'])
[ ] history2 = model2.fit(X_train, y_train,
                         epochs=5.
                         batch_size=64,
                         validation_data=(X_test, y_test))
 ⊕ Epoch 1/5
     80/80 -
                               - 15s 144ms/step - accuracy: 0.6283 - loss: 0.6456 - val_accuracy: 0.8145 - val_loss: 0.4055
     Epoch 2/5
     80/80 -
                              — 22s 161ms/step - accuracy: 0.8938 - loss: 0.2623 - val_accuracy: 0.8382 - val_loss: 0.3697
     Epoch 3/5
     80/80 -
                               - 20s 153ms/step - accuracy: 0.9656 - loss: 0.1104 - val_accuracy: 0.8540 - val_loss: 0.4207
     Epoch 4/5
     80/80 -
                              - 21s 155ms/step - accuracy: 0.9829 - loss: 0.0582 - val_accuracy: 0.8390 - val_loss: 0.4808
     Epoch 5/5
     80/80
                              - 21s 157ms/step - accuracy: 0.9933 - loss: 0.0384 - val_accuracy: 0.8421 - val_loss: 0.6642
[ ] loss, accuracy = model2.evaluate(X test, y test)
```

Model Comparison

RNN Model:

Advantages: Simple architecture that is easy to implement. Disadvantages: Struggles with capturing long-term dependencies and maintaining context over extended sequences. Best For: Suitable for tasks involving short sequences or where long-term context is not critical.

LSTM Model:

Advantages: Specifically designed to manage long-term dependencies through its gating mechanisms, which help preserve important information across longer sequences. Disadvantages: More complex and computationally intensive compared to a basic RNN. Best For: Ideal for tasks that require robust context retention and handling of longer sequences, such as textual data in fake news classification.

GRU Model:

Advantages: Offers similar performance to LSTM but with a more straightforward architecture and fewer parameters, which can lead to faster training and lower resource consumption. Disadvantages: Although efficient, its performance can vary depending on the dataset and specific task. Best For: Provides a good balance between performance and computational efficiency, making it suitable for scenarios where both performance and training efficiency are important.

Recommendations

If Performance is Critical: LSTM is generally the preferred choice due to its superior ability to handle long-term dependencies compared to basic RNNs. For datasets with long sequences that require deep context understanding, LSTM often delivers the best results.

If Training Efficiency is a Priority: GRU can be a more efficient alternative to LSTM. It is especially useful when computational resources are limited or when faster training is needed without a significant drop in performance.

For Simplicity and Baseline Comparison: Start with a basic RNN to establish a performance baseline. This approach helps to assess the improvement offered by more advanced architectures like LSTM or GRU.

Final Decision

LSTM is generally recommended for its effectiveness in managing long-term dependencies, which is crucial for comprehending context in complex text sequences. It tends to offer the best performance for tasks involving intricate textual data. Conversely, GRU provides a compelling alternative if you need a more efficient model with faster training times, offering performance similar to LSTM but with a simpler architecture.