

```
# importing dependencies
import pandas as pd
import numpy as np
from sklearn.model_selection import train_test_split
from sklearn.linear_model import LinearRegression
from sklearn.ensemble import RandomForestRegressor
from sklearn.metrics import mean_squared_error
from sklearn.preprocessing import StandardScaler
from sklearn.pipeline import Pipeline
```

```
#Load dataset
file_path = '/content/cropwatMeerut.xlsx - Sheet1.csv'
df = pd.read_csv(file_path)
```

```
print(df.head())
```

	Year	Month	Min Temp (°C)	Max Temp (°C)	Humidity (%)	Wind (km/day)	\
0	2005	January	7.7	20.1	50		6
1	2005	February	10.8	23.2	49		7
2	2005	March	17.1	30.4	45		6
3	2005	April	20.4	36.3	22		8
4	2005	May	24.8	39.5	20		9

	Sun (hours)	Rad (MJ/m <sup>2</sup> /day)	ETo (mm/day)
0	4.2	9.9	1.12
1	6.1	13.8	1.65
2	7.2	17.7	2.64
3	8.0	20.8	3.32
4	7.2	20.7	3.72

```
#checking if missing values present in dataset
print(df.isnull().sum())
```

```
Year          0
Month         0
Min Temp (°C) 0
Max Temp (°C) 0
Humidity (%)  0
Wind (km/day) 0
Sun (hours)   0
Rad (MJ/m²/day) 0
ETo (mm/day)  0
dtype: int64
```

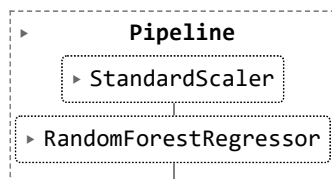
```
#dividing x(input values) parameters and Y(target) Parameters
X= df.drop(columns=['ETo (mm/day)', 'Year', 'Month'])
Y=df['ETo (mm/day)']
```

```
#Splitting data
X_Train, X_Test, Y_Train, Y_Test = train_test_split(X,Y, test_size=0.2,random_state=42)
```

```
#Creating a random forest regression model model
rf_model= RandomForestRegressor(n_estimators=100,random_state=42)
```

```
#creating pipeline for StandardScaler and RandomForestRegression
pipeline=Pipeline([
    ('scaler', StandardScaler()),
    ('rf',rf_model)
])
```

```
#training
pipeline.fit(X_Train,Y_Train)
```



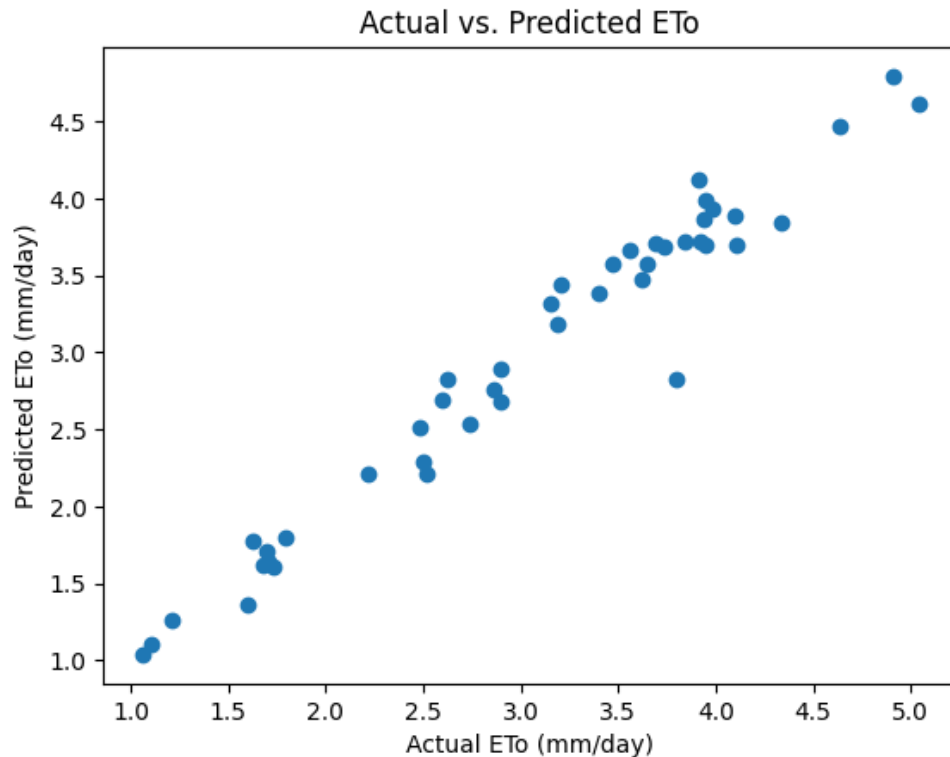
```
#making predictions
predictions=pipeline.predict(X_Test)
```

```
# Evaluate the model
mse = mean_squared_error(Y_Test, predictions)
print(f'Random Forest Regression MSE on the test set: {mse}')
```

Random Forest Regression MSE on the test set: 0.05321705522727278

```
import matplotlib.pyplot as plt

plt.scatter(Y_Test, predictions)
plt.xlabel('Actual ETo (mm/day)')
plt.ylabel('Predicted ETo (mm/day)')
plt.title('Actual vs. Predicted ETo')
plt.show()
```



## Hyperparameter Tuning

```
from sklearn.model_selection import GridSearchCV

# Define the hyperparameter grid
param_grid = {
    'rf__n_estimators': [50, 100, 150],
    'rf__max_depth': [None, 10, 20],
    'rf__min_samples_split': [2, 5, 10]
}

# Create the grid search object
grid_search = GridSearchCV(pipeline, param_grid, cv=5, scoring='neg_mean_squared_error', n_jobs=-1)

# Fit the grid search to the data
grid_search.fit(X_Train, Y_Train)

# Print the best hyperparameters
print("Best Hyperparameters:", grid_search.best_params_)

# Get the best model from grid search
best_model = grid_search.best_estimator_

# Make predictions using the best model
best_predictions = best_model.predict(X_Test)

# Evaluate the best model
best_mse = mean_squared_error(Y_Test, best_predictions)
print(f'Best Random Forest Regression MSE on the test set: {best_mse}')
```

Best Hyperparameters: {'rf\_\_max\_depth': 10, 'rf\_\_min\_samples\_split': 2, 'rf\_\_n\_estimators': 150}  
 Best Random Forest Regression MSE on the test set: 0.051938496921612495

## Let's make some predictions

```
# Assuming 'new_data' is a DataFrame containing new data with the same features as our training data
new_data = pd.DataFrame({
    'Min Temp (°C)': [7],
    'Max Temp (°C)': [20],
    'Humidity (%)': [55],
    'Wind (km/day)': [8],
    'Sun (hours)': [4],
    'Rad (MJ/m²/day)': [9.2]
})

# Using best_model to make predictions
new_predictions = best_model.predict(new_data)

# Display the predictions
print('Predicted ETo (mm/day):', new_predictions[0])

Predicted ETo (mm/day): 1.1363897924297928
```