- # Programming Language and Version: Python 3.9.6
- # Environment Setup Variables: None
- # Additional Information: The project utilizes standard Python libraries and does not require any specific environment setup variables.

Data Preprocessing

```
## Countries: China, USA, Brazil, Indonesia
## Crops: Rice, Wheat, Corn, Soya Beans
## Year: 2000-2021
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import seaborn as sns
from sklearn.impute import KNNImputer
from sklearn.preprocessing import OneHotEncoder
from sklearn.model selection import train test split
from sklearn.metrics import mean squared error, mean absolute error
from scipy.sparse import hstack, csr matrix
import warnings
# Suppress FutureWarnings
warnings.simplefilter(action='ignore', category=FutureWarning)
"""## Crop Production"""
Crop Production df = pd.read csv("Crop Production.csv")
Crop Production df.drop(['Domain Code', 'Domain', 'Area Code (M49)',
'Element Code',
                         'Item Code (CPC)', 'Year Code', 'Flag', 'Flag
Description', 'Note'], axis=1,inplace=True)
Crop Production df =
Crop Production df[Crop Production df['Year'].between(2000, 2021)]
Crop Production df.rename(columns={'Area': 'Country', 'Element':
'Prod_type', 'Item': 'Crop_Name', 'Value': 'Crop_Production_Value',
                                   'Unit': 'Crop Production Unit'},
inplace=True)
Crop Production df.head()
unique values = Crop Production df['Year'].unique()
print(unique values)
"""### *Pivoting the Prod type Column*"""
Crop Production df['Prod type'] = Crop Production df['Prod type'] + ' ' +
Crop Production df['Crop Production Unit']
```

```
# Pivot the DataFrame
Crop_Production_pivot_df = Crop Production df.pivot(index=['Country',
'Crop Name', 'Year'], columns='Prod type',
values='Crop Production Value').reset index()
# Rename columns
Crop Production pivot df.columns.name = None
# Display the pivoted DataFrame
Crop Production pivot df.head()
null values count = Crop Production pivot df.isnull().sum()
print(null values count)
# Data Types Check
print("Data Types Check:")
print(Crop Production pivot df.dtypes)
print("\n")
# Data Distribution Check
print("Data Distribution Check:")
print(Crop Production pivot df.describe())
print("\n")
## Crop Trade
Crop Trade df = pd.read csv("Crop Trade.csv")
Crop Trade df.drop(['Domain Code', 'Domain', 'Area Code (M49)', 'Element
Code',
                         'Item Code (CPC)', 'Year Code', 'Flag', 'Flag
Description', 'Note'], axis=1,inplace=True)
Crop Trade df = Crop Trade df[Crop Trade df['Year'].between(2000, 2021)]
Crop Trade df.rename(columns={'Area': 'Country', 'Element': 'Trade type',
'Item': 'Crop Name',
                              'Value': 'Trade Value', 'Unit':
'Trade Unit'}, inplace=True)
Crop Trade df.head()
unique values = Crop Trade df['Year'].unique()
print(unique values)
"""### *Pivoting the Trade type Column*"""
Crop Trade df['Trade type'] = Crop Trade df['Trade type'] + ' ' +
Crop Trade df['Trade Unit']
# Pivot the DataFrame
Crop Trade pivot df = Crop Trade df.pivot(index=['Country', 'Crop Name',
'Year'], columns='Trade type', values='Trade Value').reset index()
# Rename columns
Crop Trade pivot df.columns.name = None
```

```
# Display the pivoted DataFrame
Crop Trade pivot df.head()
# Data Types Check
print("Data Types Check:")
print(Crop Trade pivot df.dtypes)
print("\n")
null_values_count = Crop_Trade_pivot_df.isnull().sum()
print(null values count)
# Visualize distribution before handling missing values
plt.figure(figsize=(12, 6))
plt.subplot(1, 2, 1)
plt.hist(Crop Trade pivot df["Export Quantity t"], bins=20,
color='skyblue', edgecolor='black')
plt.title('Export Quantity Distribution (Before)')
plt.xlabel('Export Quantity (tonnes)')
plt.ylabel('Frequency')
plt.subplot(1, 2, 2)
plt.hist(Crop Trade pivot df["Export Value 1000 USD"], bins=20,
color='salmon', edgecolor='black')
plt.title('Export Value Distribution (Before)')
plt.xlabel('Export Value (1000 USD)')
plt.ylabel('Frequency')
plt.tight layout()
plt.show()
# Fill missing values with appropriate methods (e.g., mean, median, zero)
Crop Trade pivot df["Export
Quantity t"].fillna(Crop Trade pivot df["Export Quantity t"].mean(),
inplace=True)
Crop Trade pivot df["Export Value 1000 USD"].fillna(0, inplace=True)
Assuming missing values indicate zero export value
# Check if there are any remaining missing values
missing values = Crop Trade pivot df.isnull().sum()
print("Remaining missing values:")
print(missing values)
# Summary statistics after handling missing values
print("Summary statistics after handling missing values:")
print(Crop_Trade pivot df.describe())
print("\n")
# Visualize distribution after handling missing values
plt.figure(figsize=(12, 6))
plt.subplot(1, 2, 1)
plt.hist(Crop_Trade_pivot df["Export Quantity t"], bins=20,
color='skyblue', edgecolor='black')
plt.title('Export Quantity Distribution (After)')
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```
plt.xlabel('Export Quantity (tonnes)')
plt.ylabel('Frequency')
plt.subplot(1, 2, 2)
plt.hist(Crop Trade pivot df["Export Value 1000 USD"], bins=20,
color='salmon', edgecolor='black')
plt.title('Export Value Distribution (After)')
plt.xlabel('Export Value (1000 USD)')
plt.ylabel('Frequency')
plt.tight layout()
plt.show()
## Emission from Crops
Emission from Crops df = pd.read csv("Emission from Crops.csv")
Emission from Crops df.drop(['Domain Code', 'Domain', 'Area Code (M49)',
'Element Code', 'Source', 'Source Code',
                    'Item Code (CPC)', 'Year Code', 'Flag', 'Flag
Description', 'Note'], axis=1,inplace=True)
Emission from Crops df.rename(columns={'Area': 'Country', 'Element':
'Emission_type', 'Item': 'Crop_Name', 'Value': 'Emission_Value_kt',
'Unit': 'Emission Unit'}, inplace=True)
Emission from Crops df.head()
Emission from Crops df.drop(columns = "Emission Unit", axis = 1, inplace =
True)
Emission from Crops df.shape
Emission from Crops df.head()
null values count = Emission from Crops df.isnull().sum()
print(null values count)
# Data Types Check
print("Data Types Check:")
print(Emission from Crops df.dtypes)
print("\n")
# Summary statistics
print("Summary statistics:")
print(Emission from Crops df.describe())
print("\n")
## Land Use
Land Use df = pd.read csv("Land Use.csv")
Land Use df.drop(['Domain Code', 'Domain', 'Area Code (M49)', 'Element
Code', 'Element',
                    'Item Code', 'Year Code', 'Flag', 'Flag Description',
'Note'], axis=1, inplace=True)
```

```
Land Use df.rename(columns={'Area': 'Country','Item': 'Area type',
'Value': 'Area Value', 'Unit': 'Area Unit'}, inplace=True)
Land Use df.head()
unique values = Land Use df['Year'].unique()
print(unique values)
# Rename the "Area Value" column to "Area Value 1000 ha"
Land Use df.rename(columns={'Area Value': 'Area Value 1000 ha'},
inplace=True)
# Drop the "Area Unit" column
Land Use df.drop(columns="Area Unit", inplace=True)
Land Use df.head()
null values count = Land Use df.isnull().sum()
print(null values count)
# Data Types Check
print("Data Types Check:")
print(Land Use df.dtypes)
print("\n")
# Summary statistics
print("Summary statistics:")
print(Land Use df.describe())
print("\n")
## Pesticides Use
Pesticides Use df = pd.read csv("Pesticides Use.csv")
Pesticides Use df.drop(['Domain Code', 'Domain', 'Area Code (M49)',
'Element Code', 'Element',
                    'Item Code', 'Year Code', 'Flag', 'Flag Description',
'Note'], axis=1, inplace=True)
Pesticides Use df.rename(columns={'Area':'Country','Item':
'Pesticide Type', 'Value': 'Pesticide Value', 'Unit': 'Pesticide Unit'},
inplace=True)
Pesticides Use df.head()
unique values = Pesticides Use df['Year'].unique()
print(unique values)
# Rename the "Pesticide Value" column to "Pesticide Value t"
Pesticides Use df.rename(columns={'Pesticide Value': 'Pesticide Value t'},
inplace=True)
# Drop the "Pesticide Unit" column
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Pesticides Use df.drop(columns="Pesticide Unit", inplace=True)
Pesticides Use df.head()
null values count = Pesticides Use df.isnull().sum()
print(null values count)
# Data Types Check
print("Data Types Check:")
print(Pesticides Use df.dtypes)
print("\n")
# Summary statistics
print("Summary statistics:")
print(Pesticides Use df.describe())
print("\n")
## Value of Agricultural production
Value of Agricultural production df =
pd.read csv("Value of Agricultural Production.csv")
Value_of_Agricultural_production_df.drop(['Domain Code', 'Domain', 'Area
Code (M49)', 'Element', 'Element Code', 'Unit',
                         'Item Code (CPC)', 'Year Code', 'Flag', 'Flag
Description'], axis=1,inplace=True)
Value of Agricultural production df =
Value of Agricultural production df[Value of Agricultural production df['Y
ear'].between(2000, 2021)]
Value of Agricultural production df.rename(columns={'Area':'Country','Item
': 'Crop_Name', 'Value': 'Gross_Production_Value_1000 USD', 'Unit':
'Agri Prod Unit'}, inplace=True)
Value of Agricultural production df.head()
unique values = Value of Agricultural production df['Year'].unique()
print(unique values)
null values count = Value of Agricultural production df.isnull().sum()
print(null values count)
# Data Types Check
print("Data Types Check:")
print(Value of Agricultural production df.dtypes)
print("\n")
# Summary statistics
print("Summary statistics:")
print(Value of Agricultural production df.describe())
print("\n")
## Cropland Nutrient Balance
Cropland Nutrient Balance df =
pd.read csv("Cropland Nutrient Balance.csv")
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Cropland Nutrient Balance df.drop(['Domain Code', 'Domain', 'Area Code
(M49)', 'Element Code', 'Unit',
                         'Item Code', 'Year Code', 'Flag', 'Flag
Description', 'Note'], axis=1,inplace=True)
Cropland Nutrient Balance df =
Cropland Nutrient Balance df[Cropland Nutrient Balance df['Year'].between(
2000, 2021)]
Cropland Nutrient Balance df.rename(columns={'Area':'Country','Element':
'Cropland Nutrient Management Metrics', 'Item': 'Input Factors', 'Value':
'Fertilizer_Usage_Value_kg/ha', 'Unit': 'Agri Prod Unit'}, inplace=True)
Cropland Nutrient Balance df.head()
unique values = Cropland Nutrient Balance df['Year'].unique()
print(unique values)
null values count = Cropland Nutrient Balance df.isnull().sum()
print(null values count)
total values = Cropland Nutrient Balance df.shape[0]
percentage missing = (null values count / total values) * 100
# Display the result
print(percentage missing)
# Visualize Fertilizer Usage Distribution
plt.figure(figsize=(10, 6))
sns.boxplot(x='Input Factors', y='Fertilizer Usage Value kg/ha',
data=Cropland Nutrient Balance df)
plt.title('Distribution of Fertilizer Usage by Input Factors')
plt.xlabel('Input Factors')
plt.ylabel('Fertilizer Usage')
plt.xticks(rotation=45)
plt.show()
# Investigate Group Sizes
group sizes = Cropland Nutrient Balance df['Input Factors'].value counts()
plt.figure(figsize=(10, 6))
sns.barplot(x=group sizes.index, y=group sizes.values)
plt.title('Group Sizes by Input Factors')
plt.xlabel('Input Factors')
plt.ylabel('Group Size')
plt.xticks(rotation=45)
plt.show()
"""- The group sizes for different input factors vary significantly.
- "Mineral fertilizers" and "Manure applied to soils" have the largest
group sizes, followed by "Crop residual" and "Biological fixation."
- "Seed" and "Leaching" have relatively smaller group sizes.
11 11 11
# Visual inspection using boxplot
plt.figure(figsize=(8, 6))
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```
sns.boxplot(x=Cropland Nutrient Balance df['Fertilizer Usage Value kg/ha']
plt.title('Boxplot of Fertilizer Usage')
plt.xlabel('Fertilizer Usage')
plt.show()
"""- The boxplot reveals the presence of several outliers, with some
extreme values in the fertilizer usage data.
- The distribution of fertilizer usage appears to be right-skewed, with a
long tail towards higher values.
11 11 11
# Visualize distribution before imputation
plt.figure(figsize=(8, 6))
plt.hist(Cropland Nutrient Balance df["Fertilizer Usage Value kg/ha"].drop
na(), bins=20, color='skyblue', edgecolor='black')
plt.title('Fertilizer Usage Distribution (Before Imputation)')
plt.xlabel('Fertilizer Usage (kg/ha)')
plt.ylabel('Frequency')
plt.show()
"""### *Filling Missing Data*"""
from scipy.stats.mstats import winsorize
winsorized values =
winsorize (Cropland Nutrient Balance df['Fertilizer Usage Value kg/ha'],
                              limits=(0.05, 0.95))
# Convert the winsorized values array into a pandas Series
winsorized series = pd.Series(winsorized values)
# Fill the missing values in the original DataFrame with the winsorized
values
Cropland Nutrient Balance df['Fertilizer Usage Value kg/ha'] =
Cropland Nutrient Balance df['Fertilizer Usage Value kg/ha'].fillna(winsor
ized series)
# Verify that there are no more missing values
missing values after imputation =
Cropland Nutrient Balance df.isnull().sum()
print("Missing values after imputation:")
print(missing values after imputation)
# Visualize distribution before imputation
plt.figure(figsize=(8, 6))
plt.hist(Cropland Nutrient Balance df["Fertilizer Usage Value kg/ha"].drop
na(), bins=20, color='orange', edgecolor='black')
plt.title('Fertilizer Usage Distribution (After Imputation)')
plt.xlabel('Fertilizer Usage (kg/ha)')
plt.ylabel('Frequency')
plt.show()
```

```
merged df = pd.merge(Crop Production pivot df, Crop Trade pivot df,
on=['Country', 'Year','Crop Name'], how='inner')
merged df = pd.merge(merged df, Emission from Crops df, on=['Country',
'Year', 'Crop Name'], how='inner')
merged df = pd.merge(merged df, Value of Agricultural production df,
on=['Country', 'Year','Crop Name'], how='inner')
merged df = pd.merge(merged df, Land Use df, on=['Country', 'Year'],
how='inner')
merged df = pd.merge(merged df, Pesticides Use df, on=['Country', 'Year'],
how='inner')
merged df = pd.merge(merged df, Cropland Nutrient Balance df,
on=['Country', 'Year'], how='inner')
merged df.head()
merged df.shape
merged df.info()
"""## Feature Engineering"""
year = merged df['Year']
# Sinusoidal Transformation
merged df['Year sin'] = np.sin(2 * np.pi * year / max(year))
merged df['Year cos'] = np.cos(2 * np.pi * year / max(year))
# Display the modified DataFrame
print(merged df[['Year', 'Year sin', 'Year cos']].head())
# Emission-related Interaction Features
merged df['Emission to Production ratio'] = merged df['Emission Value kt']
/ merged df['Production t']
merged df['Emission to Area ratio'] = merged df['Emission Value kt'] /
merged df['Area harvested ha']
merged df['Emission Value per Pesticide'] = merged df['Emission Value kt']
/ merged df['Pesticide Value t']
# Land Use-related Interaction Features
merged df['Area to Production ratio'] = merged df['Area harvested ha'] /
merged df['Production t']
merged df['Area Value to Pesticide ratio'] = merged df['Area Value 1000
ha'] / merged df['Pesticide Value t']
# Pesticide-related Interaction Features
merged df['Pesticide to Production ratio'] =
merged df['Pesticide Value t'] / merged df['Production t']
# Trade related Interaction features
merged df['Export to Production ratio'] = merged df['Export Quantity t'] /
merged df['Production t']
merged df['Import to Production ratio'] = merged df['Import Quantity t'] /
merged df['Production t']
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```
# Compute aggregate statistics across different groups
# Mean production by country
country production mean =
merged df.groupby('Country')['Production t'].mean()
merged df['Country Production Mean'] =
merged df['Country'].map(country production mean)
# Median production by crop name
crop production median =
merged df.groupby('Crop Name')['Production t'].median()
merged df['Crop Production Median'] =
merged df['Crop Name'].map(crop production median)
# Sum of pesticide usage by year
yearly pesticide sum =
merged df.groupby('Year')['Pesticide Value t'].sum()
merged_df['Yearly_Pesticide_Sum'] =
merged df['Year'].map(yearly pesticide sum)
# Standard deviation of yield by country
country_yield_std = merged_df.groupby('Country')['Yield 100 g/ha'].std()
merged_df['Country Yield Std'] =
merged df['Country'].map(country yield std)
# Total emission value by crop name
crop emission total =
merged df.groupby('Crop Name')['Emission Value kt'].sum()
merged df['Crop Emission Total'] =
merged_df['Crop_Name'].map(crop_emission_total)
# Dropping duplicates
merged df.drop duplicates(inplace=True)
merged df.head()
merged df.shape
merged df.to csv('merged file.csv', index=False)
# Exploratory Data Analysis
merged df.describe()
import matplotlib.pyplot as plt
import seaborn as sns
# List of numerical columns to plot
num cols = ['Area harvested ha', 'Production t', 'Yield 100 g/ha',
            'Gross Production Value 1000 USD', 'Area Value 1000 ha',
'Pesticide Value t', 'Fertilizer Usage Value kg/ha',
            'Area_to_Production_ratio', 'Area_Value_to_Pesticide_ratio',
'Export to Production ratio',
```

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'Country Production Mean', 'Crop Production Median',
'Yearly Pesticide Sum',
            'Country Yield Std', 'Crop Emission Total']
# Setting up the matplotlib figure
plt.figure(figsize=(18, 20))
# Looping through the numerical columns to create a histogram for each
for i, col in enumerate (num cols, 1):
    plt.subplot(8, 3, i) # Adjust the grid size based on the number of
columns
    sns.histplot(merged df[col], kde=False)
    plt.title(col)
    plt.ylabel('Frequency')
    plt.xlabel(col)
plt.tight layout()
plt.show()
# Grouping by 'Crop Name' and summing the 'Production t'
total production per crop =
merged_df.groupby('Crop_Name')['Production t'].sum().reset index()
# Sorting the crops by total production in descending order
total production per crop =
total production per crop.sort values('Production t', ascending=False)
import matplotlib.pyplot as plt
import seaborn as sns
plt.figure(figsize=(8, 6))
sns.barplot(data=total production per crop, x='Crop Name',
y='Production t', color='skyblue')
plt.title('Total Production by Crop')
plt.xlabel('Total Production (tons)')
plt.ylabel('Crop Name')
plt.show()
import matplotlib.pyplot as plt
import seaborn as sns
# Group by 'Country' and 'Crop Name' and sum the production.
country crop production = merged df.groupby(['Country',
'Crop Name'])['Production t'].sum().reset index()
# Sort the production within each country and get the top N crops.
top crops per country = country crop production.groupby('Country').apply(
    lambda x: x.sort values('Production t', ascending=False).head(5) #
Taking the top 5 crops as an example
).reset index(drop=True)
# Get a sorted list of unique countries
countries = top crops per country['Country'].unique()
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# Determine the number of subplot rows we'll need, with up to 4 plots per
row for instance
n rows = len(countries) // 4 + (len(countries) % 4 > 0)
# Set up the matplotlib figure
fig, axes = plt.subplots(n rows, 4, figsize=(20, 5 * n rows)) # Adjust
the size as needed
# Flatten the axes array for easy indexing
axes = axes.flatten()
# Loop through the countries and create a plot for each
for i, country in enumerate(countries):
    # Filter the dataframe for the country
    country data = top crops per country[top crops per country['Country']
== countryl
    # Create the horizontal bar plot
    sns.barplot(x='Crop Name', y='Production t', data=country data,
ax=axes[i], palette='coolwarm')
    axes[i].set title(f'Top 5 Crops in {country}')
    axes[i].set xlabel('Crop Name')
    axes[i].set ylabel('Total Production (tons)')
# Hide any unused subplots
for j in range(i + 1, len(axes)):
    axes[j].set visible(False)
# Adjust the layout
plt.tight layout()
plt.show()
# Calculate the production per hectare for each row in the dataframe
merged df['Production per ha'] = merged df['Production t'] /
merged df['Area harvested ha']
# Group by 'Crop Name' and calculate the mean production per hectare for
each crop
productivity per crop =
merged df.groupby('Crop Name')['Production per ha'].mean().reset index()
# Sort the crops by productivity in descending order
productivity per crop sorted =
productivity per crop.sort values ('Production per ha', ascending=False)
plt.figure(figsize=(8, 5))
sns.barplot(x='Crop Name', y='Production per ha',
data=productivity per crop sorted, color='lightblue')
plt.title('Average Production (tons per hectare)')
plt.xlabel('Average Productivity per Hectare by Crop')
plt.ylabel('Crop Name')
plt.show()
# Calculate the production per hectare for each row in the dataframe
```

```
merged df['Production per ha'] = merged df['Production t'] /
merged df['Area harvested ha']
# Group by 'Country' and calculate the mean production per hectare for
each country
production per ha by country =
merged df.groupby('Country')['Production per ha'].mean().reset index()
# Sort the countries by mean production per hectare in descending order
production_per_ha_by_country_sorted =
production per ha by country.sort values('Production per ha',
ascending=False)
plt.figure(figsize=(8, 5))
sns.barplot(x='Country', y='Production per ha',
data=production per ha by country sorted.head(20), color='green')
Adjust to display the number of top countries you'd like to show
plt.title('Top Countries by Average Crop Production per Hectare')
plt.xlabel('Country')
plt.ylabel('Average Production per Hectare (tons/ha)')
plt.show()
"""## Export and Import Analysis"""
trade value = merged df.groupby('Country')[['Export Value 1000 USD',
'Import Value 1000 USD']].sum().reset index()
plt.figure(figsize=(8, 5))
# Setting the positions of the bars
barWidth = 0.4
r1 = range(len(trade value))
r2 = [x + barWidth for x in r1]
# Create the bars for exports
plt.bar(r1, trade value['Export Value 1000 USD'], color='blue',
width=barWidth, edgecolor='grey', label='Export Value')
# Create the bars for imports right next to the export bars
plt.bar(r2, trade value['Import Value 1000 USD'], color='red',
width=barWidth, edgecolor='grey', label='Import Value')
# Add labels to the x-axis at the bar centers
plt.xlabel('Country', fontweight='bold')
plt.xticks([r + barWidth / 2 for r in range(len(trade value))],
trade value['Country'], rotation=90)
plt.ylabel('Value (1000 USD)')
plt.title('Export vs Import Value for Each Country')
plt.legend()
plt.show()
# Calculate the total export and import quantity for each country
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trade balance = merged df.groupby('Country')[['Export Quantity t', 'Import
Quantity t']].sum().reset index()
# Now let's plot the grouped bar chart
plt.figure(figsize=(8, 5))
# Create the bars for exports
plt.bar(r1, trade balance['Export Quantity t'], color='blue',
width=barWidth, edgecolor='grey', label='Export Quantity')
# Create the bars for imports
plt.bar(r2, trade balance['Import Quantity t'], color='red',
width=barWidth, edgecolor='grey', label='Import Quantity')
# Add labels to the x-axis at the bar centers
plt.xlabel('Country', fontweight='bold')
plt.xticks([r + barWidth for r in range(len(trade balance))],
trade balance['Country'], rotation=90)
plt.ylabel('Quantity (tons)')
plt.yscale('log') # Apply logarithmic scale
plt.title('Export vs Import Quantity for Each Country (Logarithmic
Scale)')
# Create legend & Show graphic
plt.legend()
plt.show()
# Calculate the unit price for exports and imports
merged_df['Export_Unit_Price_USD_per_ton'] = (merged_df['Export Value_1000
USD'] * 1000) / merged df['Export Quantity t']
merged df['Import Unit Price USD per ton'] = (merged df['Import Value 1000
USD'] * 1000) / merged df['Import Quantity t']
# Handle divisions by zero or missing data if any
merged df['Export Unit Price USD per ton'].fillna(0, inplace=True)
merged df['Import Unit Price USD per ton'].fillna(0, inplace=True)
merged df.replace([np.inf, -np.inf], 0, inplace=True)
# Now we can calculate the average unit price for each country by
averaging these values
country unit prices = merged df.groupby('Country').agg({
    'Export Unit Price USD per ton': 'mean',
    'Import Unit Price USD per ton': 'mean'
}).reset index()
plt.figure(figsize=(8, 5))
# Setting the positions of the bars
barWidth = 0.35
r1 = range(len(country unit prices))
r2 = [x + barWidth for x in r1]
# Create the bars for export unit prices
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plt.bar(r1, country unit prices['Export Unit Price USD per ton'],
color='blue', width=barWidth, edgecolor='grey', label='Export Unit Price')
# Create the bars for import unit prices
plt.bar(r2, country unit prices['Import Unit Price USD per ton'],
color='red', width=barWidth, edgecolor='grey', label='Import Unit Price')
# Add labels to the x-axis at the bar centers
plt.xlabel('Country', fontweight='bold')
plt.xticks([r + barWidth for r in range(len(country unit prices))],
country unit prices['Country'], rotation=90)
plt.ylabel('Unit Price (USD/ton)')
plt.title('Average Unit Price for Export vs Import per Country')
plt.legend()
plt.show()
# Calculate the total export and import values for each crop type
crop trade values = merged df.groupby('Crop Name').agg({
    'Export Value 1000 USD': 'sum',
    'Import Value 1000 USD': 'sum'
}).reset index()
# Sort the crops by export value to see which are the highest
crop trade values.sort values('Export Value 1000 USD', ascending=False,
inplace=True)
# You could visualize the top N crops by export value to see what is
contributing to export values
plt.figure(figsize=(8, 5))
sns.barplot(x='Crop Name', y='Export Value 1000 USD',
data=crop trade values.head(20), color='blue', label='Export Value')
sns.barplot(x='Crop Name', y='Import Value 1000 USD',
data=crop trade values.head(20), color='red', alpha=0.5, label='Import
Value')
plt.xticks(rotation=90)
plt.title('Crops by Export vs Import Value')
plt.legend()
plt.show()
# Calculate the trade balance by subtracting the import value from the
export value
trade value['Trade Balance 1000USD'] = trade value['Export Value 1000
USD'] - trade value['Import Value 1000 USD']
# Now let's plot the trade balance for each country using values
plt.figure(figsize=(8, 5))
sns.barplot(x='Country', y='Trade Balance 1000USD', data=trade value,
palette='vlag')
plt.xticks(rotation=90)
plt.title('Net Trade Balance of Agricultural Products per Country
(Value)')
plt.xlabel('Country')
plt.ylabel('Trade Balance (1000 USD)')
```

```
plt.show()
# Calculate the average yield for each crop and country
average yield = merged df.groupby(['Country', 'Crop Name'])['Yield 100
g/ha'].mean().reset index()
# Sort the results by yield in descending order
average yield = average yield.sort values('Yield 100 g/ha',
ascending=False)
import seaborn as sns
import matplotlib.pyplot as plt
plt.figure(figsize=(20, 10))
sns.scatterplot(data=average yield, x='Country', y='Yield 100 g/ha',
hue='Crop Name', size='Yield 100 g/ha', sizes=(20, 200))
plt.xticks(rotation=90)
plt.title('Average Crop Yield per Country')
plt.xlabel('Country')
plt.ylabel('Yield (100g per hectare)')
plt.legend(title='Crop Name', bbox to anchor=(1.05, 1), loc='upper left')
plt.tight layout()
plt.show()
# Calculate the efficiency of pesticide usage
merged df['Pesticide Efficiency'] = merged df['Yield 100 g/ha'] /
merged df['Pesticide Value t']
# Calculate the efficiency of fertilizer usage
merged_df['Fertilizer_Efficiency'] = merged_df['Yield_100 g/ha'] /
merged df['Fertilizer Usage Value kg/ha']
# Handle potential divisions by zero or missing data
merged df.replace([np.inf, -np.inf], np.nan, inplace=True)
merged df.fillna(0, inplace=True)
# Calculate correlation matrix for yield, pesticide usage, and fertilizer
correlation matrix = merged df[['Yield 100 g/ha', 'Pesticide Value t',
'Fertilizer Usage Value kg/ha']].corr()
# Use seaborn to visualize the correlation matrix
import seaborn as sns
import matplotlib.pyplot as plt
plt.figure(figsize=(8, 4))
sns.heatmap(correlation matrix, annot=True, cmap='coolwarm', center=0)
plt.title('Correlation Matrix for Yield and Input Usage')
plt.show()
# Calculate emissions efficiency
merged df['Emissions per ton'] = merged df['Emission Value kt'] /
merged df['Production_t']
```

```
# Replace infinite values with NaN and then fill or remove them
merged df.replace([np.inf, -np.inf], np.nan, inplace=True)
merged df.dropna(subset=['Emissions per ton'], inplace=True)
# Calculate correlation matrix for production and emissions
correlation matrix prod emiss = merged df[['Production t',
'Emission Value kt']].corr()
# Visualize the correlation matrix
plt.figure(figsize=(8, 6))
sns.heatmap(correlation matrix prod emiss, annot=True, cmap='coolwarm',
plt.title('Correlation Matrix for Production and Emissions')
plt.show()
# Group data by crop or country to see average emissions per ton
average emissions per crop =
merged df.groupby('Crop Name')['Emissions per ton'].mean().reset index()
# Sort and visualize the top crops by emissions efficiency
average emissions per crop sorted =
average emissions per crop.sort values ('Emissions per ton',
ascending=False)
plt.figure(figsize=(8, 5))
sns.barplot(x='Emissions per ton', y='Crop Name',
data=average emissions per crop sorted)
plt.title('Emissions Efficiency per Crop')
plt.xlabel('Emissions per ton (kt/ton)')
plt.ylabel('Crop Name')
plt.show()
# Assuming that 'Fertilizer Usage Value kg/ha' represents the amount of
fertilizer applied per hectare
merged df['Nutrient Use Efficiency'] = merged df['Yield 100 g/ha'] /
merged_df['Fertilizer_Usage_Value_kg/ha']
# Example for a bar chart visualization
plt.figure(figsize=(8, 5))
sns.barplot(x='Crop Name', y='Nutrient Use Efficiency', data=merged df)
plt.title('Efficiency (Yield per kg/ha of Fertilizer)')
plt.xlabel('Nutrient Use Efficiency by Crop)')
plt.ylabel('Crop Name')
plt.show()
# For example, to look at production trends:
production trends =
merged df.groupby('Year')['Production t'].sum().reset index()
# For emissions trends:
emission trends =
merged df.groupby('Year')['Emission Value kt'].sum().reset index()
import matplotlib.pyplot as plt
```

```
# Line chart for production trends
plt.figure(figsize=(10, 5))
plt.plot(production trends['Year'], production trends['Production t'],
marker='o')
plt.title('Production Trends Over Time')
plt.xlabel('Year')
plt.ylabel('Total Production (tons)')
plt.grid(True)
plt.show()
# Line chart for emissions trends
plt.figure(figsize=(10, 5))
plt.plot(emission trends['Year'], emission trends['Emission Value kt'],
marker='o', color='red')
plt.title('Emission Trends Over Time')
plt.xlabel('Year')
plt.ylabel('Emissions (kt)')
plt.grid(True)
plt.show()
# Calculate the percentage of land used for each crop each year and its
impact on production.
# We will create a new feature 'Land Use Percentage' that represents the
percentage of land used for a particular crop each year.
# First, we find out the total land available each year (assuming it's the
sum of the area harvested for all crops).
total_land_each_year = merged df.groupby('Year')['Area
harvested ha'].sum().reset index().rename(columns={'Area harvested ha':
'Total Area ha'})
# Merge this total land data back into the original dataframe
merged with total land = pd.merge(merged df, total land each year,
on='Year')
# Calculate the percentage of land used for each crop each year
merged with total land['Land Use Percentage'] =
(merged with total land['Area harvested ha'] /
merged with total land['Total Area ha']) * 100
# Group by year and crop to see the impact on production
land use impact on production = merged with total land.groupby(['Year',
'Crop Name']).agg({
    'Land Use Percentage': 'mean', # Average percentage of land used for
the crop that year
    'Production t': 'sum' # Total production for the crop that
}).reset index()
# Plotting the relationship between land use percentage and production
plt.figure(figsize=(14, 7))
```

```
sns.scatterplot(data=land use impact on production,
x='Land Use Percentage', y='Production t', hue='Crop Name')
plt.title('Impact of Land Use Percentage on Crop Production')
plt.xlabel('Land Use Percentage (%)')
plt.ylabel('Total Production (tons)')
plt.legend(title='Crop Name', bbox to anchor=(1.05, 1), loc=2)
plt.grid(True)
plt.show()
# Descriptive statistics
descriptive stats = merged df.describe()
# Calculate the IQR for each numeric feature
Q1 = merged df.quantile(0.25)
Q3 = merged df.quantile(0.75)
IOR = 03 - 01
# Identify outliers for each feature
is outlier = (merged df < (Q1 - 1.5 * IQR)) | (merged df > (Q3 + 1.5 *
IQR))
import matplotlib.pyplot as plt
import seaborn as sns
# Assuming merged df is your DataFrame and it has been preprocessed
numeric features = merged df.select dtypes(include=['float64', 'int64'])
# Create violin plots for each numeric feature
for column in numeric features.columns:
    plt.figure(figsize=(8, 4))
    sns.violinplot(data=merged df, x=column)
    plt.title(f'Violin Plot for {column}')
    plt.show()
for column in is outlier:
    if (is outlier[column] == True).any():
        print(column, len(column))
# Model Building
merged df = pd.read csv("merged file.csv")
numeric cols = merged df.select dtypes(include=['float64',
'int64']).columns
# Define the percentile thresholds
lower percentile = 0.02
upper percentile = 0.98
# Capping the outliers for all numerical features in the DataFrame
for col in numeric cols:
    # Compute the 1st and 99th percentiles
    lower bound = merged df[col].quantile(lower percentile)
    upper bound = merged df[col].quantile(upper percentile)
```

```
# Cap values below the 1st percentile to the 1st percentile value
    # Cap values above the 99th percentile to the 99th percentile value
    merged df[col] = np.where(merged df[col] < lower bound, lower bound,
merged df[col])
    merged df[col] = np.where(merged df[col] > upper bound, upper bound,
merged df[col])
merged df.head()
# List of relevant features
relevant features = [
    'Country', 'Crop Name', 'Year', 'Area harvested ha', 'Production t',
'Yield 100 g/ha',
    'Export Quantity t', 'Export Value 1000 USD', 'Import Quantity t',
'Import Value 1000 USD',
    'Emission type', 'Emission Value kt', 'Gross Production Value 1000
USD', 'Area Value 1000 ha',
    'Pesticide Type', 'Pesticide Value t', 'Fertilizer Usage Value kg/ha',
    'Emission to Production ratio', 'Emission to Area ratio',
'Emission Value per Pesticide',
    'Area to Production ratio'
]
# Create a new DataFrame with only the selected features
data relevant = merged df[relevant features]
data relevant.head()
# Select only numerical features for the correlation matrix
numerical features = data relevant.select dtypes(include=['float64',
'int64']).columns
correlation matrix = data relevant[numerical features].corr()
# Plot the correlation matrix using a heatmap
plt.figure(figsize=(12, 8))
sns.heatmap(correlation matrix, annot=True, fmt=".2f", cmap='coolwarm',
cbar=True)
plt.title('Correlation Matrix of Numerical Features')
plt.show()
# Dropping highly correlated features based on heatmap analysis
columns to drop = ['Area harvested ha', 'Export Quantity t', 'Import
Value 1000 USD', 'Gross Production Value 1000 USD']
data reduced = data relevant.drop(columns=columns to drop)
# Function to calculate VIF for each feature
def calculate vif(df):
    vif data = pd.DataFrame()
    vif data['feature'] = df.columns
    vif data['VIF'] = [variance inflation factor(df.values, i) for i in
range(df.shape[1])]
    return vif data
```

```
#recalculate the VIF for the reduced dataset
vif scores reduced =
calculate vif(data reduced.select dtypes(include=['float64', 'int64',
'uint8']))
print(vif scores reduced.sort values('VIF', ascending=False))
# Dropping the specified high VIF features
columns to drop more = ['Yield 100 g/ha', 'Emission to Production ratio',
'Emission to Area ratio']
data reduced further = data reduced.drop(columns=columns to drop more)
data reduced further.head()
vif scores updated =
calculate vif(data reduced further.select dtypes(include=['float64',
'int64', 'uint8']))
print(vif scores updated.sort values('VIF', ascending=False))
# Define target variables
target vars = ['Production t', 'Export Value 1000 USD']
# Define numerical and categorical columns, excluding the target columns
numerical cols = data reduced further.select dtypes(include=['int64',
'float64']).columns.difference(target vars).tolist()
categorical cols =
data reduced further.select dtypes(include=['object']).columns.tolist()
# Create the preprocessing pipeline for numerical data
numeric transformer = Pipeline(steps=[
    ('scaler', StandardScaler()) # Standardize numerical features
])
# Create the preprocessing pipeline for categorical data
categorical transformer = Pipeline(steps=[
    ('onehot', OneHotEncoder(handle unknown='ignore')) # Encode
categorical features
])
# Combine all elements into a large transformer
preprocessor = ColumnTransformer(
   transformers=[
        ('num', numeric transformer, numerical cols),
        ('cat', categorical transformer, categorical cols)
   1)
# Split data into training and testing sets based on the year
train = data reduced further[data reduced further['Year'] <= 2019]</pre>
test = data reduced further[data reduced further['Year'] > 2019]
# Drop the 'Year' column if it's no longer needed for modeling
X_train = train.drop(['Production t', 'Export Value 1000 USD'], axis=1)
y train prod = train['Production t']
y_train_export = train['Export Value 1000 USD']
```

```
X test = test.drop(['Production t', 'Export Value 1000 USD'], axis=1)
y test prod = test['Production t']
y test export = test['Export Value 1000 USD']
"""## Linear Regression"""
# Create the full pipeline including the preprocessor and the Linear
Regression model
pipeline lr prod = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', LinearRegression())
1)
# Training the pipeline on the training data
pipeline lr prod.fit(X train, y train prod)
# Predicting on the test data
y_pred_prod = pipeline_lr_prod.predict(X_test)
# Evaluating the model
lr mae = mean absolute error(y test prod, y pred prod)
lr_mae_percentage = (lr_mae / np.mean(y_test_prod)) * 100
print("Mean Absolute Error (MAE): {:.2f}%".format(lr mae percentage))
# Mean Squared Error (MSE)
lr_mse = mean_squared_error(y_test_prod, y_pred_prod)
lr mse percentage = (lr mse / (np.mean(y test prod) ** 2)) * 100
print("Mean Squared Error (MSE): {:.2f}%".format(lr mse percentage))
# Root Mean Squared Error (RMSE)
lr rmse = np.sqrt(lr mse)
lr rmse percentage = (lr rmse / np.mean(y test prod)) * 100
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(lr rmse percentage))
r2 = r2 score(y test prod, y_pred_prod)
print("R2 Score:", r2)
"""## Lasso Regression"""
pipeline lasso = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', Lasso(alpha=1.0, random state=42))
1)
# Training the pipeline on the training data for Production
pipeline lasso.fit(X train, y train prod)
# Predicting on the test data
y pred prod = pipeline lasso.predict(X test)
# Evaluating the model
lasso mae = mean absolute error(y test prod, y pred prod)
```

```
lasso mae percentage = (lasso mae / np.mean(y test prod)) * 100
print("Mean Absolute Error (MAE): {:.2f}%".format(lasso mae percentage))
# Mean Squared Error (MSE)
lasso mse = mean squared error(y test prod, y pred prod)
lasso mse percentage = (lasso mse / (np.mean(y test prod) ** 2)) * 100
print("Mean Squared Error (MSE): {:.2f}%".format(lasso mse percentage))
# Root Mean Squared Error (RMSE)
lasso rmse = np.sqrt(lasso mse)
lasso_rmse_percentage = (lasso_rmse / np.mean(y test prod)) * 100
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(lasso rmse percentage))
r2 = r2 score(y test prod, y pred prod)
print("R2 Score:", r2)
"""## Ridge Regression"""
pipeline_ridge = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', Ridge(alpha=1.0, random state=42))
])
# Training the pipeline on the training data for Production
pipeline ridge.fit(X train, y train prod)
# Predicting on the test data
y_pred_prod = pipeline_ridge.predict(X_test)
# Evaluating the model
ridge mae = mean absolute error(y test prod, y pred prod)
ridge mae percentage = (ridge mae / np.mean(y test prod)) * 100
print("Mean Absolute Error (MAE): {:.2f}%".format(ridge mae percentage))
# Mean Squared Error (MSE)
ridge_mse = mean_squared_error(y_test_prod, y_pred_prod)
ridge mse percentage = (ridge mse / (np.mean(y test prod) ** 2)) * 100
print("Mean Squared Error (MSE): {:.2f}%".format(ridge mse percentage))
# Root Mean Squared Error (RMSE)
ridge rmse = np.sqrt(ridge mse)
ridge rmse percentage = (ridge rmse / np.mean(y test prod)) * 100
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(ridge rmse percentage))
r2 = r2_score(y_test_prod, y pred prod)
print("R2 Score:", r2)
"""## SVM Regressor"""
pipeline svr = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', SVR(kernel='poly'))
```

```
])
# Training the pipeline on the training data for Production
pipeline_svr.fit(X_train, y_train_prod)
# Predicting on the test data
y pred prod = pipeline svr.predict(X test)
# Evaluating the model
svr mae = mean absolute error(y test prod, y pred prod)
svr mae percentage = (svr mae / np.mean(y test prod)) * 100
print("Mean Absolute Error (MAE): {:.2f}%".format(svr mae percentage))
# Mean Squared Error (MSE)
svr_mse = mean_squared_error(y_test_prod, y_pred_prod)
svr_mse_percentage = (svr_mse / (np.mean(y_test_prod) ** 2)) * 100
print("Mean Squared Error (MSE): {:.2f}%".format(svr_mse_percentage))
# Root Mean Squared Error (RMSE)
svr rmse = np.sqrt(svr mse)
svr_rmse_percentage = (svr_rmse / np.mean(y_test_prod)) * 100
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(svr_rmse_percentage))
r2 = r2 score(y test prod, y pred prod)
print("R2 Score:", r2)
"""## XGBoost Regressor"""
pipeline xgb = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', XGBRegressor(objective='reg:squarederror',
random state=42))
])
# Training the pipeline on the training data for Production
pipeline xgb.fit(X train, y train prod)
# Predicting on the test data
y pred prod = pipeline xgb.predict(X test)
xgb_mae = mean_absolute_error(y_test_prod, y_pred_prod)
xgb mae percentage = (xgb mae / np.mean(y test prod)) * 100
print("Mean Absolute Error (MAE): {:.2f}%".format(xgb_mae_percentage))
# Mean Squared Error (MSE)
xgb_mse = mean_squared_error(y_test_prod, y_pred_prod)
xgb mse percentage = (xgb mse / (np.mean(y test prod) ** 2)) * 100
print("Mean Squared Error (MSE): {:.2f}%".format(xgb mse percentage))
# Root Mean Squared Error (RMSE)
xgb rmse = np.sqrt(xgb mse)
xgb rmse percentage = (xgb rmse / np.mean(y test prod)) * 100
```

```
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(xgb rmse percentage))
r2 = r2 score(y test prod, y pred prod)
print("R2 Score:", r2)
# Define the parameter grid
param grid = {
    'regressor n estimators': [100, 200],
    'regressor__max_depth': [3, 5, 7],
    'regressor learning rate': [0.01, 0.1],
    'regressor subsample': [0.7, 0.9],
    'regressor colsample bytree': [0.7, 0.9]
}
# Setup the grid search
grid search = GridSearchCV(pipeline xgb, param grid, cv=3,
scoring='neg mean squared error')
grid search.fit(X train, y train prod)
print("Best parameters:", grid search.best params )
print("Best cross-validation score: ", np.sqrt(-grid search.best score ))
# Update pipeline with the best parameters
pipeline xgb optimized = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', XGBRegressor(
        objective='reg:squarederror',
        colsample bytree=0.9,
        learning_rate=0.1,
        \max depth=5,
        n estimators=200,
        subsample=0.9,
        random state=42))
1)
# Training the optimized pipeline on the training data
pipeline xgb_optimized.fit(X_train, y_train_prod)
# Predicting on the test data
y pred prod = pipeline xgb optimized.predict(X test)
xgb optimized mae = mean absolute error(y test prod, y pred prod)
xgb optimized mae percentage = (xgb optimized mae / np.mean(y test prod))
* 100
print("Mean Absolute Error (MAE):
{:.2f}%".format(xgb optimized mae percentage))
# Mean Squared Error (MSE)
xgb optimized mse = mean squared error(y test prod, y pred prod)
xqb optimized mse percentage = (xqb optimized mse / (np.mean(y test prod)
** 2)) * 100
print("Mean Squared Error (MSE):
{:.2f}%".format(xgb optimized mse percentage))
```

```
# Root Mean Squared Error (RMSE)
xgb optimized rmse = np.sqrt(xgb optimized mse)
xgb optimized rmse percentage = (xgb optimized rmse /
np.mean(y test prod)) * 100
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(xgb optimized rmse percentage))
r2 = r2 score(y test prod, y pred prod)
print("R2 Score:", r2)
"""## Random Forest"""
pipeline rf = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', RandomForestRegressor(n estimators=100, random state=42,
max depth= 6))
1)
# Training the pipeline on the training data for Production
pipeline rf.fit(X train, y train prod)
# Predicting on the test data
y pred prod = pipeline rf.predict(X test)
# Evaluating the model
rf mae = mean absolute error(y test prod, y pred prod)
rf mae percentage = (rf mae / np.mean(y test prod)) * 100
print("Mean Absolute Error (MAE): {:.2f}%".format(rf_mae_percentage))
# Mean Squared Error (MSE)
rf mse = mean squared error(y test prod, y pred prod)
rf mse percentage = (rf mse / (np.mean(y test prod) ** 2)) * 100
print("Mean Squared Error (MSE): {:.2f}%".format(rf mse percentage))
# Root Mean Squared Error (RMSE)
rf rmse = np.sqrt(rf mse)
rf rmse percentage = (rf rmse / np.mean(y test prod)) * 100
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(rf rmse percentage))
r2 = r2 score(y test prod, y pred prod)
print("R2 Score:", r2)
# Calculate residuals
residuals = y test prod - y pred prod
# Plotting the histogram of residuals
plt.figure(figsize=(6, 4))
plt.hist(residuals, bins=30, alpha=0.5, color='g')
plt.title('Histogram of Prediction Errors')
plt.xlabel('Prediction Error')
plt.ylabel('Frequency')
```

```
plt.show()
# Boxplot of residuals
plt.figure(figsize=(6, 4))
plt.boxplot(residuals, vert=False)
plt.title('Boxplot of Prediction Errors')
plt.xlabel('Prediction Error')
plt.show()
plt.figure(figsize=(10, 6))
stats.probplot(residuals, dist="norm", plot=plt)
plt.title('Q-Q Plot')
plt.show()
plt.figure(figsize=(10, 6))
plt.scatter(y pred prod, residuals)
plt.title('Residuals vs. Predicted Values')
plt.xlabel('Predicted Values')
plt.ylabel('Residuals')
plt.axhline(y=0, color='r', linestyle='--')
plt.show()
df analysis = pd.DataFrame({
  'predicted_values': y_pred_prod,
  'residuals': residuals
# Define quantile-based ranges for predicted values
df analysis['predicted range'] = pd.qcut(df analysis['predicted values'],
4)
# Calculate summary statistics for each range
summary stats =
df analysis.groupby('predicted range')['residuals'].agg(['mean', 'median',
'std', 'count'])
# Plot the distribution of residuals for each range
fig, axes = plt.subplots(nrows=4, figsize=(8, 12))
for (range val, subset), ax in zip(df analysis.groupby('predicted range'),
axes):
    ax.hist(subset['residuals'], bins=10, alpha=0.7)
    ax.set title(f'Residuals for Predicted Range: {range val}')
    ax.set xlabel('Residuals')
    ax.set ylabel('Frequency')
plt.tight layout()
plt.show()
# Identifying high error instances
high error threshold = residuals.abs().mean() + 1.5 *
residuals.abs().std() # Using mean + 1.5*STD as threshold
high error indices = residuals[residuals.abs() >
high error threshold].index
high error data = X test.loc[high error indices]
```

```
# Analyzing features of high error instances
feature analysis = high error data.describe(include='all')
display(feature analysis)
def calculate metrics(y t, y p):
    model mae = mean absolute error(y t, y p)
    model mae percentage = (model mae / np.mean(y t)) * 100
    # Mean Squared Error (MSE)
    model mse = mean squared error(y t, y p)
    model mse percentage = (model mse / (np.mean(y t) ** 2)) * 100
    # Root Mean Squared Error (RMSE)
    model rmse = np.sqrt(model mse)
    model rmse percentage = (model rmse / np.mean(y t)) * 100
    model_r2 = r2_score(y_t, y_p)
    return model mae percentage, model mse percentage,
model rmse percentage, model r2
"""## Residual Calculation and Grouping"""
df test = []
df test = pd.DataFrame({
    'actual': y test prod,
    'predicted': y_pred_prod
})
df test = pd.concat([X test, df test], axis = 1)
df test.shape
df test['residuals'] = df test['actual'] - df test['predicted']
# Group residuals into bins (for example, by quartiles)
df test['residual group'] = pd.qcut(df test['residuals'], 4, labels=['Q1',
'\overline{2}', '\overline{3}', '\overline{4}'])
# Plotting the distribution of residuals within each bin
plt.figure(figsize=(12, 6))
sns.boxplot(x='residual group', y='residuals', data=df test)
plt.title('Distribution of Residuals by Group')
plt.xlabel('Residual Group')
plt.ylabel('Residuals')
plt.grid(True)
plt.show()
"""## Extract Data Points from Q1 and Q4:"""
# Filter data for Q1 and Q4 groups
df q1 = df test[df test['residual group'] == 'Q1']
df q4 = df test[df test['residual group'] == 'Q4']
# Descriptive statistics for Q1
```

```
print("Descriptive Statistics for Q1:")
display(df q1.describe())
# Descriptive statistics for Q4
print("Descriptive Statistics for Q4:")
display(df q4.describe())
# Correlation matrix for the Q1 group
corr q1 = df test[df test['residual group'] == 'Q1'].corr()
print("Correlation Matrix for Q1:")
display(corr q1)
# Correlation matrix for the Q4 group
corr q4 = df test[df test['residual group'] == 'Q4'].corr()
print("Correlation Matrix for Q4:")
display(corr q4)
# Heatmap visualization for Q1
plt.figure(figsize=(8, 6))
sns.heatmap(corr q1.drop(['predicted', 'residuals'],
axis=0).drop(['actual', 'predicted', 'residuals'], axis=1),
            annot=True,
            cmap='coolwarm', fmt=".2f")
plt.title('Correlation Heatmap for Q1')
plt.show()
# Heatmap visualization for Q4
plt.figure(figsize=(10, 8))
sns.heatmap(corr q4.drop(['predicted', 'residuals'],
axis=0).drop(['actual', 'predicted', 'residuals'], axis=1),
            annot=True, cmap='coolwarm', fmt=".2f")
plt.title('Correlation Heatmap for Q4')
plt.show()
y pred train = pipeline rf.predict(X train)
mae rf train, mse rf train, rmse rf train, r2 rf train =
calculate metrics(y train prod, y pred train)
print(f"Segment 1 - MAE: {mae rf train}, MSE: {mse rf train}, RMSE:
{rmse rf train}, R2: {r2 rf train}")
residual train = y train prod - y pred train
df test train = pd.DataFrame({
    'actual': y train_prod,
    'predicted': y pred train
})
df test train = pd.concat([X train, df test train], axis = 1)
df test train.shape
```

```
df_test_train['residuals'] = df_test_train['actual'] -
df_test_train['predicted']

# Group residuals into bins (for example, by quartiles)
df_test_train['residual_group'] = pd.qcut(df_test_train['residuals'], 2,
labels=['Q1', 'Q2'])

# Plotting the distribution of residuals within each bin
plt.figure(figsize=(6, 5))
sns.boxplot(x='residual_group', y='residuals', data=df_test_train)
plt.title('Distribution of Residuals by Group')
plt.xlabel('Residual Group')
plt.ylabel('Residuals')
plt.grid(True)
plt.show()
```

Model Optimization

```
** ** **
Based on the EDA, performing the following actions
merged df = pd.read csv("merged file.csv")
# Assuming 'merged df' is your preprocessed DataFrame and you have already
identified numeric cols
numeric cols = merged df.select dtypes(include=['float64',
'int64']).columns
# Define the percentile thresholds
lower percentile = 0.02
upper percentile = 0.98
# Capping the outliers for all numerical features in the DataFrame
for col in numeric cols:
    \# Compute the \overline{1}st and 99th percentiles
    lower bound = merged df[col].quantile(lower percentile)
    upper bound = merged df[col].quantile(upper percentile)
    # Cap values below the 1st percentile to the 1st percentile value
    # Cap values above the 99th percentile to the 99th percentile value
    merged df[col] = np.where(merged df[col] < lower bound, lower bound,
merged df[col])
    merged df[col] = np.where(merged df[col] > upper bound, upper bound,
merged df[col])
merged df.head()
# List of relevant features
relevant features = [
    'Country', 'Crop Name', 'Year', 'Area harvested ha', 'Production t',
'Yield 100 g/ha',
```

```
'Export Quantity t', 'Export Value 1000 USD', 'Import Quantity t',
'Import Value 1000 USD',
    'Emission type', 'Emission Value kt', 'Gross Production Value 1000
USD', 'Area Value 1000 ha',
    'Pesticide Type', 'Pesticide Value t', 'Fertilizer Usage Value kg/ha',
    'Emission to Production ratio', 'Emission to Area ratio',
'Emission Value per Pesticide',
    'Area to Production ratio'
]
# Create a new DataFrame with only the selected features
data relevant = merged df[relevant features]
data relevant.head()
# Select only numerical features for the correlation matrix
numerical features = data relevant.select dtypes(include=['float64',
'int64']).columns
correlation matrix = data relevant[numerical features].corr()
# Plot the correlation matrix using a heatmap
plt.figure(figsize=(12, 8))
sns.heatmap(correlation_matrix, annot=True, fmt=".2f", cmap='coolwarm',
cbar=True)
plt.title('Correlation Matrix of Numerical Features')
plt.show()
# Dropping highly correlated features based on heatmap analysis
columns to drop = ['Area harvested ha', 'Export Quantity t', 'Import
Value_1000 USD', 'Gross_Production_Value_1000 USD']
data reduced = data relevant.drop(columns=columns to drop)
# Function to calculate VIF for each feature
def calculate vif(df):
    vif data = pd.DataFrame()
    vif data['feature'] = df.columns
    vif data['VIF'] = [variance inflation factor(df.values, i) for i in
range(df.shape[1])]
    return vif data
#recalculate the VIF for the reduced dataset
vif scores reduced =
calculate_vif(data_reduced.select dtypes(include=['float64', 'int64',
'uint8']))
print(vif scores reduced.sort values('VIF', ascending=False))
# Dropping the specified high VIF features
columns to drop more = ['Yield 100 g/ha', 'Emission to Production ratio',
'Emission to Area ratio']
data reduced further = data reduced.drop(columns=columns to drop more)
data reduced further.head()
```

```
vif scores updated =
calculate vif(data reduced further.select dtypes(include=['float64',
'int64', 'uint8']))
print(vif scores updated.sort values('VIF', ascending=False))
# Define target variables
target vars = ['Production t']
# Define numerical and categorical columns, excluding the target columns
numerical_cols = data_reduced further.select dtypes(include=['int64',
'float64']).columns.difference(target vars).tolist()
categorical cols =
data reduced further.select dtypes(include=['object']).columns.tolist()
# Create the preprocessing pipeline for numerical data
numeric transformer = Pipeline(steps=[
    ('scaler', StandardScaler()) # Standardize numerical features
1)
# Create the preprocessing pipeline for categorical data
categorical transformer = Pipeline(steps=[
    ('onehot', OneHotEncoder(handle unknown='ignore')) # Encode
categorical features
1)
# Combine all elements into a large transformer
preprocessor = ColumnTransformer(
    transformers=[
        ('num', numeric transformer, numerical cols),
        ('cat', categorical transformer, categorical cols)
    ])
# Split data into training and testing sets based on the year
train = data reduced further[data reduced further['Year'] <= 2019]</pre>
test = data reduced further[data reduced further['Year'] > 2019]
# Drop the 'Year' column if it's no longer needed for modeling
X train = train.drop(['Production t'], axis=1)
y train = train['Production t']
X test = test.drop(['Production t'], axis=1)
y test = test['Production t']
"""## Random Forest"""
def calculate metrics(y t, y p):
    model mae = mean absolute error(y t, y p)
    model mae percentage = (model_mae / np.mean(y_t)) * 100
    # Mean Squared Error (MSE)
    model mse = mean squared error(y t, y p)
    model mse percentage = (model mse / (np.mean(y t) ** 2)) * 100
    # Root Mean Squared Error (RMSE)
```

```
model_rmse = np.sqrt(model mse)
    model rmse percentage = (model rmse / np.mean(y t)) * 100
    model r2 = r2 score(y t, y p)
    return model mae percentage, model mse percentage,
model rmse percentage, model r2
pipeline rf = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', RandomForestRegressor(n estimators=100, random state=42,
max depth=8))
])
# Training the pipeline on the training data for Production
pipeline rf.fit(X train, y train)
# Predicting on the train data
y pred train = pipeline rf.predict(X train)
rf mae train, rf mse train, rf rmse train, rf r2 train =
calculate metrics(y train, y pred train)
print("Mean Absolute Error (MAE): {:.2f}%".format(rf_mae_train))
print("Mean Squared Error (MSE): {:.2f}%".format(rf mse train))
print("Root Mean Squared Error (RMSE): {:.2f}%".format(rf rmse train))
print("R2 Score:", rf r2 train)
y pred test = pipeline rf.predict(X test)
rf mae test, rf mse test, rf rmse test, rf r2 test =
calculate_metrics(y test, y pred test)
print("Mean Absolute Error (MAE): {:.2f}%".format(rf mae test))
print("Mean Squared Error (MSE): {:.2f}%".format(rf mse test))
print("Root Mean Squared Error (RMSE): {:.2f}%".format(rf rmse test))
print("R2 Score:", rf_r2_test)
# Boxplot for target variable
plt.figure(figsize=(4, 4))
sns.boxplot(data reduced further['Production t'])
plt.title('Box Plot of Target Variable')
plt.show()
# Scatter plot for target variable against another feature that you
suspect might be influencing outliers
plt.figure(figsize=(8, 4))
plt.scatter(data reduced further['Emission Value kt'],
data reduced further['Production t'])
plt.xlabel('Feature')
plt.ylabel('Target')
plt.title('Scatter Plot of Feature vs. Target')
plt.show()
```

```
# Calculate IQR for the target variable
Q1 = data reduced further['Production t'].quantile(0.05)
Q3 = data reduced further['Production t'].quantile(0.95)
IQR = Q3 - Q1
# Define thresholds for outliers
lower bound = Q1 - 1.5 * IQR
upper bound = Q3 + 1.5 * IQR
# Filtering the data
filtered df = data reduced further[(data reduced further['Production t']
>= lower bound) & (data reduced further['Production t'] <= upper bound)]
print(f"Original data points: {len(data reduced further)}")
print(f"Data points after outlier removal: {len(filtered df)}")
# Split data into training and testing sets based on the year
train filter = filtered df[filtered df['Year'] <= 2019]</pre>
test filter = filtered df[filtered df['Year'] > 2019]
X train filter = train filter.drop(['Production t'], axis=1)
y train filter = train filter['Production t']
X test filter = test filter.drop(['Production t'], axis=1)
y test filter = test filter['Production t']
y pred train filter = pipeline rf.predict(X train filter)
rf mae train filter, rf mse train filter, rf rmse train filter,
rf r2 train filter = calculate metrics(y train filter,
y pred train filter)
print("Mean Absolute Error (MAE): {:.2f}%".format(rf mae train filter))
print("Mean Squared Error (MSE): {:.2f}%".format(rf mse train filter))
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(rf rmse train filter))
print("R2 Score:", rf r2 train filter)
y pred test filter = pipeline rf.predict(X test filter)
rf mae test filter, rf mse test filter, rf rmse test filter,
rf_r2_test_filter = calculate_metrics(y_test_filter, y_pred_test_filter)
print("Mean Absolute Error (MAE): {:.2f}%".format(rf mae test filter))
print("Mean Squared Error (MSE): {:.2f}%".format(rf mse test filter))
print("Root Mean Squared Error (RMSE):
{:.2f}%".format(rf rmse test filter))
print("R2 Score:", rf r2 test filter)
import matplotlib.pyplot as plt
plt.hist(filtered df['Production t'], bins=30)
plt.title('Histogram of Target Variable')
plt.show()
```

```
# Define thresholds
threshold = 1.5e8
# Create subsets
def divide segments (threshold, segment no):
    segment1 = filtered df[filtered df['Production t'] <= threshold]</pre>
    segment2 = filtered df[filtered df['Production t'] > threshold] #&
(filtered df['Production t'] <= threshold2)]</pre>
# segment3 = filtered df[(filtered df['Production_t'] > threshold)
    if segment no == 1:
        return segment1
    else:
        return segment2
X1 = divide segments(threshold,1).drop('Production t', axis=1)
y1 = divide segments(threshold,1)['Production t']
X2 = divide segments(threshold,2).drop('Production t', axis=1)
y2 = divide segments(threshold,2)['Production t']
pipeline lr = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', LinearRegression())
])
pipeline lasso = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', Lasso(alpha=1.0, random state=42))
])
# pipeline for segment 2: Ridge Regressor
pipeline ridge = Pipeline(steps = [
    ('preprocessor', preprocessor),
    ('regressor', Ridge(alpha=1.0, random state=42))
1)
# Pipeline for segment 2: Random Forest Regressor
pipeline rf = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', RandomForestRegressor(n estimators=100, random state=42,
max depth= 6))
1)
# Pipeline for segment 3: Gradient Boosting Regressor
pipeline gb = Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', GradientBoostingRegressor(n estimators=100,
random state=42))
1)
# Function to split data and train pipeline
def train test pipeline (X, y, pipeline, mode):
    X train, X test, y train, y test = train test split(X, y,
test size=0.3, random state=42)
    pipeline.fit(X train, y train)
```

```
y pred tr = pipeline.predict(X train)
   y pred = pipeline.predict(X test)
   mae tr, mse tr, rmse tr, r2 tr = calculate metrics(y train, y pred tr)
   mae, mse, rmse, r2 = calculate metrics(y test, y pred)
   if mode == 'test':
        return mae, mse, rmse, r2
    if mode == 'train':
       return mae tr, mse tr, rmse tr, r2 tr
# Applying pipelines
mae1 tr, mse1 tr, rmse1 tr, r21 tr = train test pipeline(X1, y1,
pipeline gb, 'train')
mae2 tr, mse2 tr, rmse2 tr, r22_tr = train_test_pipeline(X2, y2,
pipeline ridge, 'train')
\# mae3 tr, mse2 tr, rmse2_tr, r23_tr =
train test pipeline(segment3.drop('Production t', axis=1),
segment3['Production t'], pipeline gb)
print(f"Segment 1 - MAE: {mae1 tr}, MSE: {mse1 tr}, RMSE: {rmse1 tr}, R2:
{r21 tr}")
print(f"Segment 2 - MAE: {mae2 tr}, MSE: {mse2 tr}, RMSE: {rmse2 tr}, R2:
{r22 tr}")
# print(f"Segment 3 - MAE: {mae2 tr}, MSE: {mse2 tr}, RMSE: {rmse2 tr},
R2: {r22 tr}")
# Applying pipelines
mae1, mse1, rmse1, r21 = train test pipeline(X1,y1, pipeline gb, 'test')
mae2, mse2, rmse2, r22 = train test pipeline(X2, y2, pipeline ridge,
'test')
\# mae3, mse2, rmse2, r23 =
train test pipeline(segment3.drop('Production t', axis=1),
segment3['Production t'], pipeline gb)
print(f"Segment 1 - MAE: {mae1}, MSE: {mse1}, RMSE: {rmse1}, R2: {r21}")
print(f"Segment 2 - MAE: {mae2}, MSE: {mse2}, RMSE: {rmse2}, R2: {r22}")
# print(f"Segment 3 - MAE: {mae2}, MSE: {mse2}, RMSE: {rmse2}, R2: {r22}")
"""## Finding Threshold based on Clustering"""
from sklearn.cluster import DBSCAN
def find density based threshold(data):
   values = data['Production t'].values.reshape(-1, 1)
   clustering = DBSCAN(eps=0.5, min samples=10).fit(values)
    labels = clustering.labels
    # Find transition points between clusters
   unique labels = np.unique(labels)
    core samples = np.array([values[labels == label].mean() for label in
unique labels if label != -1])
    if len(core samples) > 1:
        sorted samples = np.sort(core samples)
        # Threshold between the first two clusters
        threshold = np.mean(sorted samples[:2])
        return threshold
```

```
thrshold db = find density based threshold(filtered df)
thrshold db
X db1 = divide segments(thrshold db,1).drop('Production t', axis=1)
y db1 = divide segments(thrshold db,1)['Production t']
X db2 = divide segments(thrshold db,2).drop('Production t', axis=1)
y db2 = divide segments(thrshold db,2)['Production t']
# Applying pipelines
mae1 db tr, mse1 db tr, rmse1 db tr, r21 db tr =
train test pipeline(X_db1, y_db1,pipeline_rf, 'train')
mae2 db tr, mse2 db tr, rmse2 db tr, r22 db tr =
train test pipeline(X db2, y db2, pipeline ridge, 'train')
\# mae3 tr, mse2 tr, rmse2 tr, r23 tr =
train test pipeline(segment3.drop('Production t', axis=1),
segment3['Production t'], pipeline gb)
print(f"Segment 1 - MAE: {mae1 db tr}, MSE: {mse1 db tr}, RMSE:
{rmse1 db tr}, R2: {r21 db tr}")
print(f"Segment 2 - MAE: {mae2 db tr}, MSE: {mse2 db tr}, RMSE:
{rmse2 db tr}, R2: {r22 db tr}")
# print(f"Segment 3 - MAE: {mae2 tr}, MSE: {mse2 tr}, RMSE: {rmse2 tr},
R2: {r22 tr}")
# Applying pipelines
mae1 db, mse1 db, rmse1 db, r21 db = train test pipeline(X db1,
y db1,pipeline gb, 'test')
mae2 db, mse2 db, rmse2 db, r22 db = train test pipeline(X db2, y db2,
pipeline ridge, 'test')
\# mae3 tr, mse2 tr, rmse2 tr, r23 tr =
train test pipeline(segment3.drop('Production t', axis=1),
segment3['Production t'], pipeline gb)
print(f"Segment 1 - MAE: {mae1 db}, MSE: {mse1 db}, RMSE: {rmse1 db}, R2:
{r21 db}")
print(f"Segment 2 - MAE: {mae2 db}, MSE: {mse2 db}, RMSE: {rmse2 db}, R2:
{r22 db}")
# print(f"Segment 3 - MAE: {mae2 tr}, MSE: {mse2 tr}, RMSE: {rmse2 tr},
R2: {r22 tr}")
"""## Change Point Detection"""
pip install ruptures
from ruptures import Pelt
import ruptures as rpt
def find change point threshold(data):
    values = data['Production t'].values
    algo = rpt.Pelt(model="12").fit(values)
    result = algo.predict(pen=10)
    # Find the first significant change point as the threshold
    if result:
        threshold = values[result[0]]
        return threshold
```

```
thr cpd = find change point threshold(filtered df)
thr_cpd
X cpd1 = divide segments(thr cpd,1).drop('Production t', axis=1)
y cpd1 = divide segments(thr cpd,1)['Production t']
X cpd2 = divide segments(thr cpd,2).drop('Production t', axis=1)
y cpd2 = divide segments(thr cpd,2)['Production t']
# Applying pipelines
mae1 cpd tr, mse1 cpd tr, rmse1 cpd tr, r21 cpd tr =
train test pipeline(X cpd1, y cpd1, pipeline rf, 'train')
mae2 cpd tr, mse2 cpd tr, rmse2 cpd tr, r22 cpd tr =
train test pipeline(X cpd2, y cpd2, pipeline rf, 'train')
# mae3_tr, mse2_tr, rmse2_tr, r23_tr =
train test pipeline(segment3.drop('Production t', axis=1),
segment3['Production_t'], pipeline_gb)
print(f"Segment 1 - MAE: {mae1 cpd tr}, MSE: {mse1 cpd tr}, RMSE:
{rmse1 cpd tr}, R2: {r21 cpd tr}")
print(f"Segment 2 - MAE: {mae2_cpd_tr}, MSE: {mse2_cpd_tr}, RMSE:
{rmse2 cpd tr}, R2: {r22 cpd tr}")
# print(f"Segment 3 - MAE: {mae2 tr}, MSE: {mse2 tr}, RMSE: {rmse2 tr},
R2: {r22 tr}")
# Applying pipelines
mae1 cpd, mse1 cpd, rmse1 cpd, r21 cpd = train test pipeline(X cpd1,
y_cpd1,pipeline_rf, 'train')
mae2 cpd, mse2 cpd, rmse2 cpd, r22 cpd = train test pipeline(X cpd2,
y cpd2, pipeline ridge, 'train')
\# mae3 tr, mse2 tr, rmse2 tr, r23 tr =
train test pipeline(segment3.drop('Production t', axis=1),
segment3['Production t'], pipeline gb)
print(f"Segment 1 - MAE: {mae1 cpd}, MSE: {mse1 cpd}, RMSE: {rmse1 cpd},
R2: {r21 cpd}")
print(f"Segment 2 - MAE: {mae2 cpd}, MSE: {mse2 cpd}, RMSE: {rmse2 cpd},
R2: {r22 cpd}")
# print(f"Segment 3 - MAE: {mae2 tr}, MSE: {mse2 tr}, RMSE: {rmse2 tr},
R2: {r22 tr}")
"""### Segregating Train and Test based on Years"""
data b 2019 = filtered df[filtered df["Year"] < 2019]</pre>
data a 2019 = filtered df[filtered df["Year"] >= 2019]
import matplotlib.pyplot as plt
plt.hist(data b 2019['Production t'], bins=30)
plt.title('Histogram of Target Variable')
plt.show()
import matplotlib.pyplot as plt
plt.hist(data a 2019['Production_t'], bins=30)
```

```
plt.title('Histogram of Target Variable')
plt.show()
thr b 2019 = find change point threshold(data b 2019)
thr b 2019
thr a 2019 = find change point threshold(data a 2019)
thr a 2019
def train and evaluate(X, y, pipeline):
    pipeline.fit(X, y)
    y pred = pipeline.predict(X)
    return calculate metrics (y, y pred)
def evaluate models(data, pipelines):
    X = data.drop(columns=['Production t'])
    y = data['Production t']
    results = {}
    for name, pipeline in pipelines.items():
        metrics = train and evaluate(X, y, pipeline)
        results[name] = metrics
    return results
# Segmenting data based on thresholds
data b 2019 segment1 = data b 2019[data b 2019["Production t"] <=
thr b 2019]
data b 2019 segment2 = data b 2019[data b 2019["Production t"] >
thr b 2019]
data a 2019 segment1 = data a 2019[data a 2019["Production t"] <=
thr_a 2019]
data a 2019 segment2 = data a 2019[data a 2019["Production t"] >
thr a 2019]
pipelines = {
    'Linear Regression': Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', LinearRegression())
]),
    'Lasso': Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', Lasso(alpha=1.0, random state=42))
]),
    'Ridge': Pipeline(steps = [
    ('preprocessor', preprocessor),
    ('regressor', Ridge(alpha=1.0, random state=42))
]),
    'Random Forest': Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', RandomForestRegressor(n estimators=100, random state=42,
max depth= 4))
```

```
]),
    'Gradient Boosting': Pipeline(steps=[
    ('preprocessor', preprocessor),
    ('regressor', GradientBoostingRegressor(n estimators=100,
random state=42))
])
}
# Evaluate models for each segment
results b segment1 = evaluate models(data b 2019 segment1, pipelines)
results b segment2 = evaluate models(data b 2019 segment2, pipelines)
results a segment1 = evaluate models(data a 2019 segment1, pipelines)
results a segment2 = evaluate models(data a 2019 segment2, pipelines)
# Print results
print("Results for data before 2019, Segment 1:", results b segment1)
print("Results for data before 2019, Segment 2:", results_b_segment2)
print("Results for data after 2019, Segment 1:", results a segment1)
print("Results for data after 2019, Segment 2:", results a segment2)
def find best model(results):
   best model name = None
   best model score = -np.inf # Initialize with negative infinity for
comparison
   best model metrics = None
    for model name, metrics in results.items():
        r2 score = metrics[-1] # The last metric in the tuple is R^2
        if r2 score > best model score:
           best model score = r2 score
            best model name = model name
            best model metrics = metrics
    return best model name, best model metrics
# Print the best model for each segment
def print best model results(segment name, results):
   best model name, best model metrics = find best model(results)
   print(f"Best model for {segment name}:")
   print(f"Model: {best model name}")
   print(f"Metrics (MAE%, MSE%, RMSE%, R²): {best model metrics}\n")
print best model results ("data before 2019, Segment 1",
results b segment1)
print best model results ("data before 2019, Segment 2",
results b segment2)
print best model results("data after 2019, Segment 1", results a segment1)
print best model results("data after 2019, Segment 2", results a segment2)
```