Problem Statement

The objective of this project is to build a predictive model that determines how likely an individual is to be a user of **illicit drugs** based on their demographic attributes, personality traits, and behavioral characteristics. The study leverages a dataset containing information on drug usage patterns (e.g., cannabis, cocaine, heroin, etc.), demographic variables (age, gender, education, etc.), and psychometric scores (e.g., neuroticism, extraversion, sensation-seeking).

This project focuses on answering the following key questions:

- 1. Which features—demographics, personality traits, or behavioral metrics—are the strongest predictors of illicit drug use?
- 2. How well can a supervised learning model predict the likelihood of illicit drug use, and what are its key performance metrics?

Scope of Illicit Drugs

"Illicit drugs" in this analysis refer to substances such as cannabis, cocaine, heroin, LSD, ketamine, methamphetamines, and similar drugs, excluding legal substances like alcohol and nicotine.

Significance

Substance abuse is a significant public health issue, with illicit drug usage often leading to adverse social, psychological, and economic consequences. Identifying predictors of illicit drug use can help inform targeted interventions, educational programs, and public health policies to prevent and reduce substance abuse.

This predictive model will provide insights into the attributes associated with illicit drug use and offer a tool to predict such behaviors, which could be utilized in further research or by public health organizations.

Project Goals

- 1. Perform **exploratory data analysis (EDA)** to uncover patterns and relationships in the data, particularly focusing on predictors of illicit drug use.
- 2. Build and evaluate a classification model to predict illicit drug use (binary classification: user vs. non-user).
- 3. Identify and interpret the most influential features contributing to predictions, using techniques like feature importance analysis or SHAP (SHapley Additive Explanations).
- 4. Address issues like **class imbalance** using techniques such as SMOTE, undersampling, or class-weighted models.

Deliverables

- 1. A cleaned and preprocessed dataset, with binary labels indicating illicit drug use.
- 2. Visualizations and analysis summarizing the relationships between predictors and drug use (e.g., correlation matrices, pair plots, bar charts).
- 3. A machine learning model (e.g., logistic regression, random forest, or XGBoost) that achieves acceptable performance based on evaluation metrics (e.g., precision, recall, F1 score, ROC-AUC).
- 4. A discussion of insights derived from the analysis and model, including the implications for public health.

In [1]:

```
from ucimlrepo import fetch_ucirepo
import pandas as pd
import seaborn as sns
from matplotlib import pyplot as plt
```

```
from sklearn.model_selection import train_test_split
from sklearn.linear_model import LogisticRegression
from sklearn.metrics import classification_report, roc_auc_score, confusion_matrix, roc_c
urve, auc

# Fetch dataset
drug_consumption_quantified = fetch_ucirepo(id=373)

# Load data as Pandas DataFrames
X = drug_consumption_quantified.data.features
y = drug_consumption_quantified.data.targets

# Merge features and targets into a single DataFrame
df = pd.concat([X, y], axis=1)
```

```
In [2]:
```

```
# Slight cleaning
df = df.rename(columns={'impuslive': 'impulsive'})
df = df.drop(columns=['semer'])
```

Load and Preview the Data

Understand the dataset structure by summarizing it

```
In [3]:
```

```
# Check the structure of the dataset
print(df.info())

# Check for missing values
print(df.isnull().sum())

# Preview the dataset
print(df.head())
```

<class 'pandas.core.frame.DataFrame'>
RangeIndex: 1885 entries, 0 to 1884
Data columns (total 30 columns):

1 gender 1885 non-null float64 2 education 1885 non-null float64 3 country 1885 non-null float64 4 ethnicity 1885 non-null float64 5 nscore 1885 non-null float64 6 escore 1885 non-null float64 7 oscore 1885 non-null float64 8 ascore 1885 non-null float64 9 cscore 1885 non-null float64 10 impulsive 1885 non-null float64	aca	COTAMINS (CC	Juan .	JO COLUMNIS,	
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2		meth	1885	non-null	object
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	20	nicotino	1005	non-null	abiaat

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                               object
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memory usage: 441.9+ KB
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             0
gender
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            0
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nscore
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cscore
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impulsive
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             0
alcohol
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amphet
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             0
cannabis
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choc
            0
coke
crack
ecstasy
heroin
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ketamine
             0
legalh
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meth
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mushrooms
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nicotine
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2 \quad 0.49788 \quad -0.48246 \quad -0.05921 \quad 0.96082 \quad -0.31685 \quad -0.46725 \quad 0.80523 \quad -0.84732
3 -0.95197 0.48246 1.16365 0.96082
                                         -0.31685 -0.14882 -0.80615 -0.01928
                     1.98437 0.96082
                                         -0.31685 0.73545 -1.63340 -0.45174
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1 0.76096 -0.14277 ...
                          CL0
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                                                            CLO CL2 CL3
2 -1.62090 -1.01450 ...
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3 0.59042 0.58489 ...
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4 -0.30172 1.30612
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                 CLO CLO
3
       CL0
                 CL2
                     CL0
       CL2
                CL2 CL0
[5 rows x 30 columns]
In [4]:
df.columns
Out[4]:
'amphet', 'amyl', 'benzos', 'caff', 'cannabis', 'choc', 'coke', 'crack', 'ecstasy', 'heroin', 'ketamine', 'legalh', 'lsd', 'meth', 'mushrooms',
       'nicotine', 'vsa'],
```

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dtype='object')

object

Understand the Target Variable

Before diving into predictors, analyze the distribution of the target variable (y).

```
In [5]:
```

```
# Define a list of drugs (update based on your analysis)
drugs = ['cannabis', 'coke', 'crack', 'heroin', 'meth', 'ecstasy',
        'lsd', 'amyl', 'ketamine', 'mushrooms', 'vsa']
# Use at least once in last year
drug user = ['CL3', 'CL4', 'CL5', 'CL6']
# Define a binary target variable for any drug use in the last year
def check drug use(row):
   for user in drug user: # Iterate over the drug user codes
       if user in row.values: # Check if the code is in the row's values
           return 1 # Return 1 if a match is found
    return 0 # Return 0 if no matches are found
# Apply the function only to the specified drug columns
df['drug use'] = df[drugs].apply(check drug use, axis=1)
dfDrugUse = df.drop(columns={'alcohol',
       'amphet', 'amyl', 'benzos', 'caff', 'cannabis', 'choc', 'coke', 'crack',
       'ecstasy', 'heroin', 'ketamine', 'legalh', 'lsd', 'meth', 'mushrooms',
       'nicotine', 'vsa'})
dfDrugUse.head()
```

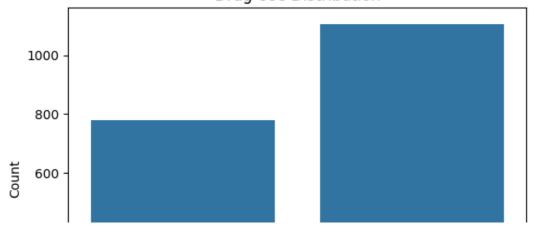
Out[5]:

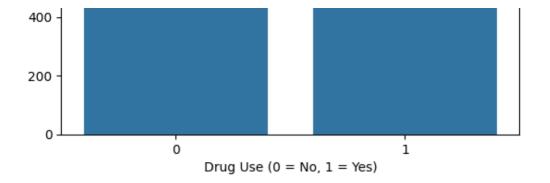
	age	gender	education	country	ethnicity	nscore	escore	oscore	ascore	cscore	impulsive	SS	drug_use
(0.49788	0.48246	-0.05921	0.96082	0.12600	0.31287	- 0.57545	- 0.58331	- 0.91699	0.00665	-0.21712	- 1.18084	0
1	0.07854	- 0.48246	1.98437	0.96082	-0.31685	- 0.67825	1.93886	1.43533	0.76096	- 0.14277	-0.71126	- 0.21575	1
2	0.49788	- 0.48246	-0.05921	0.96082	-0.31685	- 0.46725	0.80523	- 0.84732	- 1.62090	- 1.01450	-1.37983	0.40148	1
:	0.95197	0.48246	1.16365	0.96082	-0.31685	- 0.14882	0.80615	0.01928	0.59042	0.58489	-1.37983	- 1.18084	0
4	0.49788	0.48246	1.98437	0.96082	-0.31685	0.73545	- 1.63340	- 0.45174	0.30172	1.30612	-0.21712	- 0.21575	1

In [6]:

```
# Plot the new binary target variable
sns.countplot(x=df['drug_use'], y=None, hue=None, data=df)
plt.title("Drug Use Distribution")
plt.xlabel("Drug Use (0 = No, 1 = Yes)")
plt.ylabel("Count")
plt.show()
```

Drug Use Distribution





Overview of the Target Variable

• Start by checking the distribution of the binary target variable to ensure no severe class imbalance.

```
In [7]:
```

```
# Check class distribution
print(df['drug_use'].value_counts(normalize=True))

drug_use
1    0.587268
0    0.412732
Name: proportion, dtype: float64
```

Mapping Fields for Better Interpretability

In [8]:

```
# Mapping age values to age groups
age mapping = {
    -0.95197: "18-24",
   -0.07854: "25-34",
    0.49788: "35-44",
    1.09449: "45-54",
    1.82213: "55-64",
     2.59171: "65+"
}
gender mapping = {
   0.48246: "Female",
   -0.48246: "Male"
education mapping = {
    -2.43591: "Left before 16",
    -1.73790: "Left at 16",
    -1.43719: "Left at 17",
    -1.22751: "Left at 18",
    -0.61113: "Some college",
    -0.05921: "Professional cert",
    0.45468: "University degree",
    1.16365: "Masters degree",
    1.98437: "Doctorate"
country mapping = {
   -0.09765: "Australia",
    0.24923: "Canada",
   -0.46841: "New Zealand",
    -0.28519: "Other",
    0.21128: "Ireland",
    0.96082: "UK",
    -0.57009: "USA"
ethnicity mapping = {
```

```
-0.50212: "Asian",

-1.10702: "Black",

1.90725: "Mixed-Black/Asian",

0.12600: "Mixed-White/Asian",

-0.22166: "Mixed-White/Black",

0.11440: "Other",

-0.31685: "White"
```

Create a Mapped DataFrame for Visualizations

Copy the original df to a new DataFrame, dfMapped, and apply the mappings only to this new DataFrame.

```
In [9]:
```

```
# Create a separate DataFrame for mapped values
dfMapped = df.copy()

# Apply mappings to the copied DataFrame for visualizations
dfMapped['age'] = dfMapped['age'].map(age_mapping)
dfMapped['gender'] = dfMapped['gender'].map(gender_mapping)
dfMapped['education'] = dfMapped['education'].map(education_mapping)
dfMapped['country'] = dfMapped['country'].map(country_mapping)
dfMapped['ethnicity'] = dfMapped['ethnicity'].map(ethnicity_mapping)
```

Feature Distribution by Illicit Drug Use

• Analyze the distribution of demographic and personality traits with respect to illicit drug use.

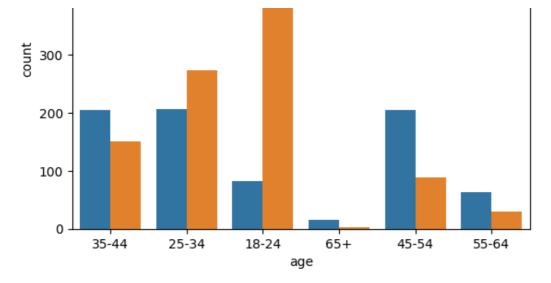
a. Demographics

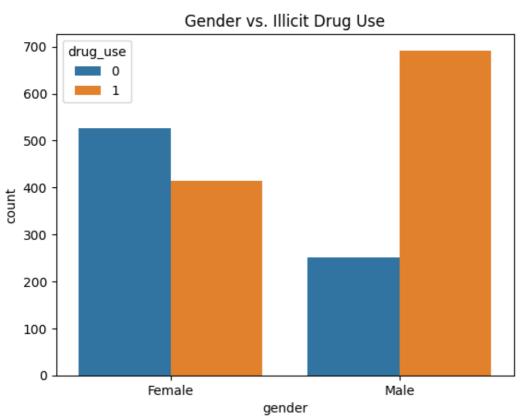
```
In [10]:
```

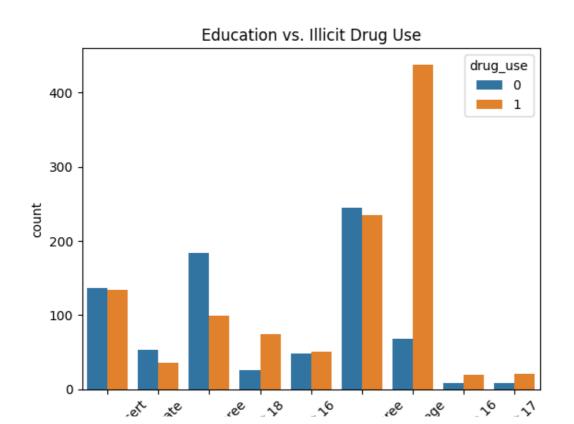
```
import seaborn as sns
import matplotlib.pyplot as plt
# Visualize age distribution by illicit drug use
sns.countplot(data=dfMapped, x='age', hue=df['drug use'])
plt.title("Age vs. Illicit Drug Use")
plt.show()
# Visualize gender distribution by illicit drug use
sns.countplot(data=dfMapped, x='gender', hue=df['drug use'])
plt.title("Gender vs. Illicit Drug Use")
plt.show()
# Visualize education distribution by illicit drug use
sns.countplot(data=dfMapped, x='education', hue=df['drug use'])
plt.title("Education vs. Illicit Drug Use")
plt.xticks(rotation=45)
plt.show()
# Visualize ethnicity distribution by illicit drug use
sns.countplot(data=dfMapped, x='ethnicity', hue=df['drug use'])
plt.title("Ethnicity vs. Illicit Drug Use")
plt.xticks(rotation=45)
plt.show()
```

Age vs. Illicit Drug Use

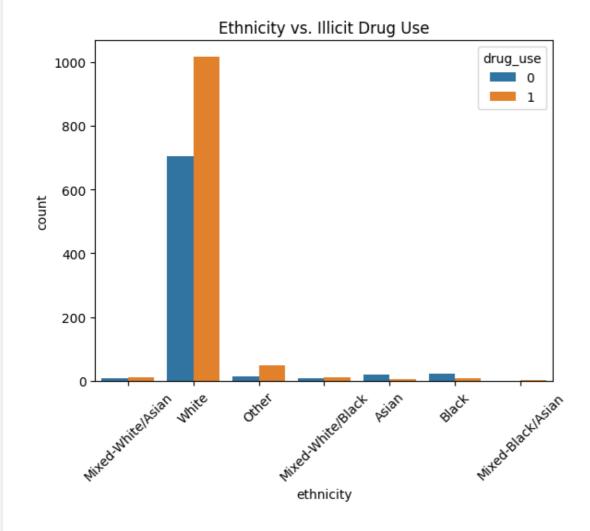












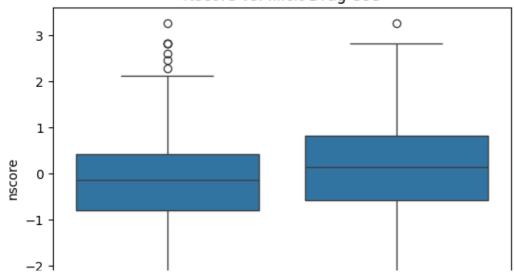
b. Personality Traits

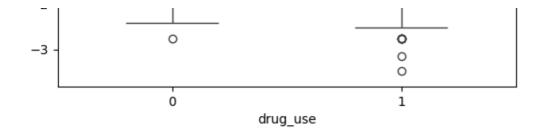
In [11]:

```
# Box plots for personality traits by illicit drug use
traits = ['nscore', 'escore', 'ascore', 'cscore', 'impulsive', 'ss']

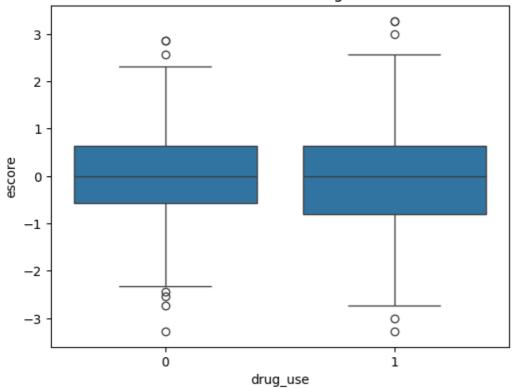
for trait in traits:
    sns.boxplot(data=df, x=df['drug_use'], y=trait)
    plt.title(f"{trait.capitalize()} vs. Illicit Drug Use")
    plt.show()
```

Nscore vs. Illicit Drug Use

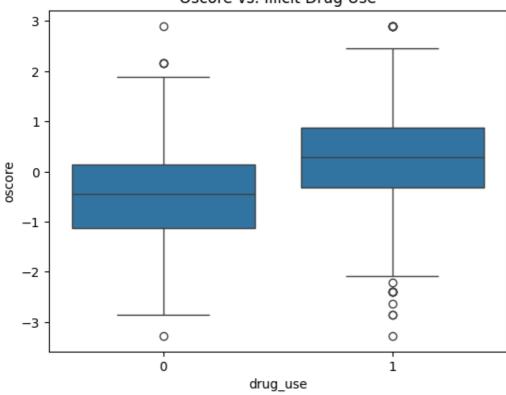




Escore vs. Illicit Drug Use

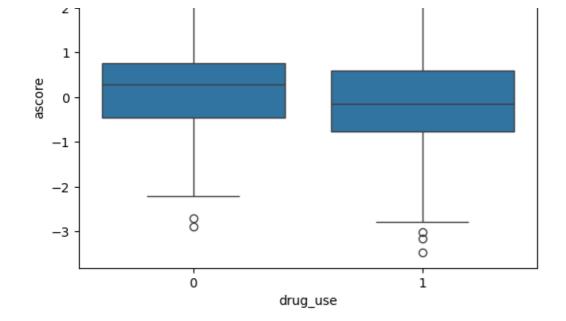


Oscore vs. Illicit Drug Use

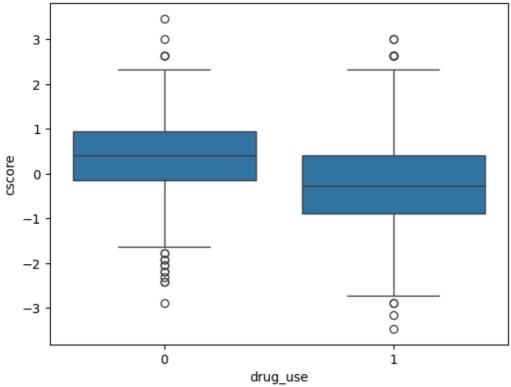


Ascore vs. Illicit Drug Use

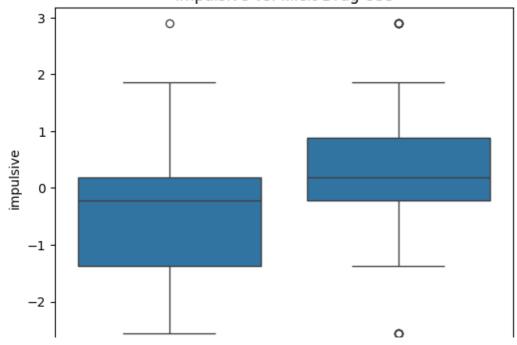


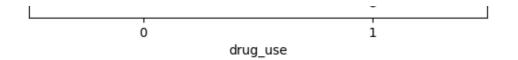


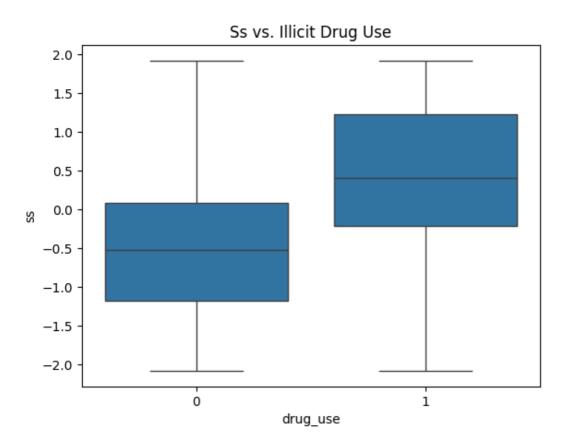




Impulsive vs. Illicit Drug Use







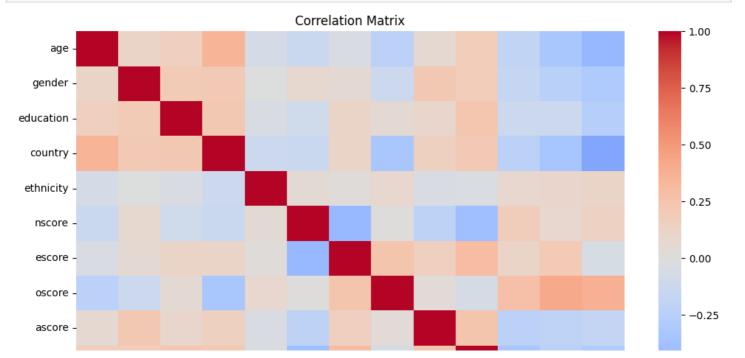
Correlation Analysis

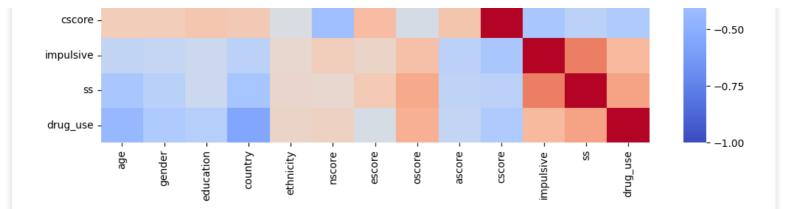
• Analyze the correlation between numerical features and illicit drug use to identify strong predictors.

```
In [12]:
```

```
# Compute correlation matrix
correlation_matrix = dfDrugUse.corr()

# Plot heatmap for correlations
plt.figure(figsize=(12, 8))
sns.heatmap(correlation_matrix, annot=False, cmap='coolwarm', vmin=-1, vmax=1)
plt.title("Correlation_Matrix")
plt.show()
```



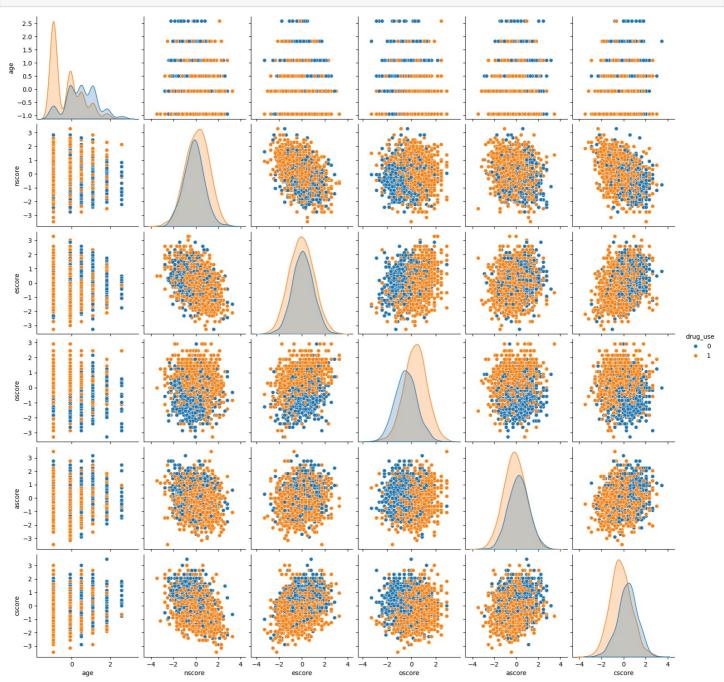


Pairwise Relationships

• Use pair plots to explore how key features relate to illicit drug use.

In [13]:

```
# Pair plot for selected features
selected_features = ['age', 'nscore', 'escore', 'oscore', 'ascore', 'cscore', 'drug_use']
sns.pairplot(df[selected_features], hue='drug_use', diag_kind='kde')
plt.show()
```



Relationships Between Features and illicit drug use:

```
In [14]:
```

```
for col in ['age', 'gender', 'education']:
    print(f"Distribution of {col} by illicit_drug_use:")
    print(dfMapped.groupby([col, 'drug_use']).size().unstack())
```

```
Distribution of age by illicit drug use:
drug_use 0
                1
age
18-24
25-34
207 2.
35-44
205 151
45-54
205 89
63 30
55-64 63
65+ 16
                2
Distribution of gender by illicit drug use:
drug_use 0 1
gender
Female 527 415
Male
         251 692
Distribution of education by illicit drug use:
                 0
education
                   53 36
Doctorate
Left at 16
                   48 51
Left at 17
Left at 18
                    9 21
                   26 74
Left before 16 9
Masters degree 184
                    9 19
                         99
Professional cert 136 134
                    68 438
Some college
University degree 245 235
```

Logistic Regression

- Type: Linear Model
- Best For:
 - Simple relationships between predictors and the target variable.
 - Interpretable models with clear feature importance (via coefficients).
 - Problems with linearly separable data.
- Pros:
 - Easy to implement and interpret.
 - Computationally efficient and fast to train.
 - Performs well with simple or moderately complex datasets.
 - Naturally outputs probabilities for classification.
- Cons:
 - Assumes linearity between predictors and the target.
 - Sensitive to multicollinearity among features.
 - May struggle with non-linear decision boundaries.

In [15]:

```
# Define features (X) and target (y)
X = dfDrugUse.drop('drug_use', axis=1)  # Features
y = dfDrugUse['drug_use']  # Target

# Split the data 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, stratify=y)
```

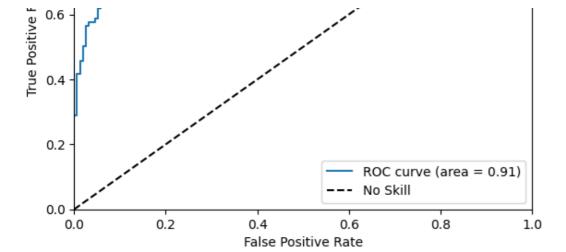
In [16]:

```
# Initialize Logistic Regression model
logistic_model = LogisticRegression(solver='liblinear', max_iter=1000, random_state=42,
class_weight='balanced')
```

```
# Train the model
logistic model.fit(X train, y train)
Out[16]:
                              LogisticRegression
LogisticRegression(class weight='balanced', max iter=1000, random state=42,
                   solver='liblinear')
In [17]:
# Predict on the test set
y pred = logistic model.predict(X test)
y_proba = logistic_model.predict_proba(X_test)[:, 1] # Probability estimates for ROC-AUC
# Classification report
print("Classification Report:")
print(classification report(y test, y pred))
# Confusion matrix
conf matrix = confusion matrix(y test, y pred)
print("Confusion Matrix:")
print(conf matrix)
Classification Report:
             precision
                          recall f1-score
                                               support
           0
                   0.76
                             0.84
                                       0.80
                                                   156
                             0.81
                                       0.84
                                                   221
           1
                   0.88
   accuracy
                                       0.82
                                                   377
  macro avg
                   0.82
                             0.82
                                       0.82
                                                   377
weighted avg
                   0.83
                             0.82
                                       0.82
                                                   377
Confusion Matrix:
[[131 25]
 [ 42 179]]
In [18]:
# Predict probabilities on the test set
y pred proba = logistic model.predict proba(X test)[:, 1]
# Calculate ROC curve
fpr, tpr, thresholds = roc curve(y test, y pred proba)
roc auc = auc(fpr, tpr)
# Plot the ROC curve
plt.figure()
plt.plot(fpr, tpr, label='ROC curve (area = %0.2f)' % roc auc)
plt.plot([0, 1], [0, 1], 'k--', label='No Skill')
plt.xlim([0.0, 1.0])
plt.ylim([0.0, 1.05])
plt.xlabel('False Positive Rate')
plt.ylabel('True Positive Rate')
plt.title('ROC Curve for Drug Use Prediction')
plt.legend()
plt.show()
                   ROC Curve for Drug Use Prediction
```

1.0

0.8



In [19]:

```
from sklearn.model_selection import GridSearchCV

param_grid = {'C': [0.01, 0.1, 1, 10], 'penalty': ['12'], 'solver': ['lbfgs']}
grid = GridSearchCV(LogisticRegression(max_iter=1000, class_weight='balanced'), param_grid, scoring='roc_auc', cv=5)
grid.fit(X_train, y_train)
print(f"Best Parameters: {grid.best_params_}")
print(f"Best ROC-AUC Score: {grid.best_score_}")

Best Parameters: {'C': 10, 'penalty': '12', 'solver': 'lbfgs'}
Best ROC-AUC Score: 0.9136074623778778
```

In [20]:

```
best_logistic_model = LogisticRegression(C=10, penalty='12', solver='lbfgs')
best_logistic_model.fit(X_train, y_train)
# Predict on the test set
y_pred = best_logistic_model.predict(X_test)
y_proba = best_logistic_model.predict_proba(X_test)[:, 1] # Probability estimates for RO
C-AUC

# Classification report
print("Classification Report:")
print(classification_report(y_test, y_pred))
# Confusion matrix
conf_matrix = confusion_matrix(y_test, y_pred)
print("Confusion Matrix:")
print(conf_matrix)
```

Classification Report:

	precision	recall	fl-score	support
0 1	0.80	0.81	0.81 0.86	156 221
accuracy macro avg weighted avg	0.83	0.83	0.84 0.83 0.84	377 377 377

Confusion Matrix:
[[126 30]
 [31 190]]

Random Forest

- Type: Ensemble Model (Bagging)
- Best For:
 - Complex, non-linear relationships.
 - Handling missing data and unbalanced datasets.
 - Feature importance analysis.

- Pros:
 - Can capture non-linear patterns and interactions between features.
 - Robust to overfitting with enough trees.
 - Automatically handles feature scaling and normalization.
 - Resistant to outliers and noise.
- Cons:
 - Can be slower to train for large datasets.
 - Interpretability is lower compared to simpler models like Logistic Regression.
 - May require hyperparameter tuning (e.g., number of trees, max depth) for optimal performance.

In [21]:

Out[21]:

RandomForestClassifier

i

RandomForestClassifier(class weight='balanced', random state=42)

In [22]:

```
# Predict on the test set
y_pred = rf_model.predict(X_test)
y_proba = rf_model.predict_proba(X_test)[:, 1] # Probability estimates for ROC-AUC

# Classification report
print("Classification_Report:")
print(classification_report(y_test, y_pred))

# ROC-AUC score
roc_auc = roc_auc_score(y_test, y_proba)
print(f"ROC-AUC Score: {roc_auc}")

# Confusion matrix
conf_matrix = confusion_matrix(y_test, y_pred)
print("Confusion Matrix:")
print(conf_matrix)
```

Classification Report:

support	f1-score	recall	precision	
156	0.82	0.79	0.84	0
221	0.88	0.89	0.86	1
377	0.85			accuracy
377 377	0.85 0.85	0.84 0.85	0.85 0.85	macro avg weighted avg

```
ROC-AUC Score: 0.9092992226476388
```

Confusion Matrix:

[[124 32]

```
[ 24 17/]]
In [23]:
from sklearn.model selection import GridSearchCV
# Define hyperparameter grid
param grid = {
    'n estimators': [100, 200, 300],
    'max_depth': [None, 10, 20, 30],
    'min_samples_split': [2, 5, 10],
    'min_samples_leaf': [1, 2, 4]
# Grid search
grid search = GridSearchCV(
   estimator=RandomForestClassifier(random state=42, class weight='balanced'),
   param grid=param grid,
   scoring='roc auc',
   cv=5,
    verbose=1,
    n jobs=2
grid search.fit(X train, y train)
# Best parameters and score
print(f"Best Parameters: {grid search.best params }")
print(f"Best ROC-AUC Score: {grid search.best score }")
Fitting 5 folds for each of 108 candidates, totalling 540 fits
Best Parameters: {'max_depth': 10, 'min_samples_leaf': 2, 'min_samples_split': 10, 'n_est
imators': 200}
Best ROC-AUC Score: 0.9102986413238237
In [24]:
best rf model = RandomForestClassifier(
   n estimators=200, # Number of trees
   random state=42,
    min samples leaf=2, # Minimum samples in leaf nodes
    min samples split=10, # Minimum samples to split nodes
    class weight='balanced', # Handle class imbalance
    max depth=10,
# Train the model
best_rf_model.fit(X_train, y_train)
Out[24]:
                      RandomForestClassifier
RandomForestClassifier(class weight='balanced', max depth=10,
                       min samples leaf=2, min samples split=10,
                       n estimators=200, random state=42)
In [25]:
# Predict on the test set
y pred = best rf model.predict(X test)
y proba = best rf model.predict proba(X test)[:, 1] # Probability estimates for ROC-AUC
# Classification report
print("Classification Report:")
```

```
# Predict on the test set
y_pred = best_rf_model.predict(X_test)
y_proba = best_rf_model.predict_proba(X_test)[:, 1] # Probability estimates for ROC-AUC

# Classification report
print("Classification Report:")
print(classification_report(y_test, y_pred))

# ROC-AUC score
roc_auc = roc_auc_score(y_test, y_proba)
print(f"ROC-AUC Score: {roc_auc}")
```

```
conf matrix = confusion matrix(y test, y pred)
print("Confusion Matrix:")
print(conf_matrix)
Classification Report:
                         recall f1-score
             precision
                                             support
                 0.81 0.83
0.88 0.86
                                  0.82
           0
                                                 156
                                      0.87
                                                 221
                                      0.85
                                                 377
   accuracy
                  0.84 0.85
                                      0.84
  macro avg
                                                 377
                            0.85
                                      0.85
                                                 377
weighted avg
                  0.85
ROC-AUC Score: 0.9150713539853811
Confusion Matrix:
[[130 26]
 [ 31 190]]
```

Support Vector Machine (Soft Margin)

• Type: Kernel-based Model

Confusion matrix

- Best For:
 - High-dimensional data.
 - Non-linear decision boundaries (via kernel functions).
 - Problems where a soft margin (allowing some misclassifications) is beneficial.
- Pros:
 - Effective for high-dimensional spaces and small datasets.
 - Can model complex relationships using kernels (e.g., radial basis function, polynomial).
 - Well-suited for imbalanced datasets with proper tuning (e.g., class weights).
- Cons:
 - Computationally expensive for large datasets.
 - Sensitive to the choice of hyperparameters (e.g., kernel type, C parameter).
 - Less interpretable compared to Logistic Regression or Random Forest.

```
In [26]:
```

```
import numpy as np
from sklearn.svm import SVC
from sklearn.model_selection import cross_val_score, GridSearchCV, train_test_split
from sklearn.metrics import classification_report, confusion_matrix, roc_auc_score
import matplotlib.pyplot as plt
```

In [27]:

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42
, stratify=y)

# Train SVM with default hyperparameters
svm_model = SVC(C=1.0, kernel='rbf', gamma='scale', probability=True, random_state=42)
svm_model.fit(X_train, y_train)

# Cross-validation
scores = cross_val_score(svm_model, X_train, y_train, cv=5)
print("Cross-validation mean accuracy: {:.3f}".format(np.mean(scores)))
```

Cross-validation mean accuracy: 0.836

In [28]:

```
# Predict on the test set
y_pred = svm_model.predict(X_test)
y_proba = svm_model.decision_function(X_test)

# Classification report
print("Classification Report:")
```

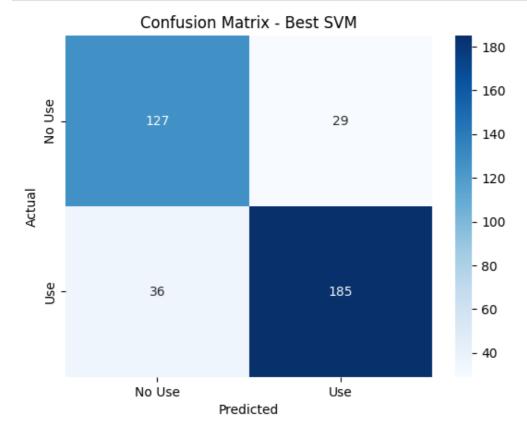
```
print(classification_report(y_test, y_pred))
# Confusion matrix
conf matrix = confusion matrix(y test, y pred)
print("Confusion Matrix:")
print(conf matrix)
# ROC-AUC Score
roc auc = roc auc score(y test, y proba)
print(f"ROC-AUC Score: {roc auc}")
Classification Report:
             precision
                         recall f1-score
                                             support
                          0.79
                   0.81
                                     0.80
                                                  156
           1
                   0.86
                           0.86
                                     0.86
                                                  221
                                      0.84
                                                  377
   accuracy
   macro avg
                  0.83
                           0.83
                                     0.83
                                                  377
                  0.84
                           0.84
                                     0.84
                                                 377
weighted avg
Confusion Matrix:
[[124 32]
 [ 30 191]]
ROC-AUC Score: 0.8974649031210117
In [29]:
kernels = ['linear', 'poly', 'rbf', 'sigmoid']
for kernel in kernels:
    svm = SVC(C=1.0, kernel=kernel, gamma='scale', random state=42)
    scores = cross val score(svm, X train, y train, cv=5)
    print(f"kernel = {kernel}, cross-val mean-accuracy: {np.mean(scores):.3f}")
kernel = linear, cross-val mean-accuracy: 0.839
kernel = poly, cross-val mean-accuracy: 0.810
kernel = rbf, cross-val mean-accuracy: 0.836
kernel = sigmoid, cross-val mean-accuracy: 0.796
In [30]:
# Define the parameter grid
param grid = {
    'C': np.logspace(-5, 5, num=11, base=2), # C values: 2^-5 to 2^5
    'gamma': np.logspace(-5, 5, num=11, base=2)  # gamma values: 2^-5 to 2^5
# GridSearchCV for hyperparameter tuning
grid = GridSearchCV(
    estimator=SVC(kernel='linear', probability=True, random state=42),
    param grid=param grid,
   cv=5,
    scoring='roc auc',
    verbose=1
# Fit the grid search to the data
grid.fit(X_train, y_train)
# Print the best parameters and corresponding accuracy
print("Best parameters found: ", grid.best params )
print("Best cross-validation ROC-AUC: {:.3f}".format(grid.best score ))
# Use the best model for predictions
best model = grid.best estimator
y pred best = best model.predict(X test)
print("Test set accuracy of the best model: {:.3f}".format(best model.score(X test, y tes
t)))
Fitting 5 folds for each of 121 candidates, totalling 605 fits
```

Best parameters found: {'C': 2.0, 'gamma': 0.03125}

```
Best cross-validation ROC-AUC: 0.914
Test set accuracy of the best model: 0.828
```

In [31]:

```
# Confusion matrix for the best model
conf_matrix_best = confusion_matrix(y_test, y_pred_best)
sns.heatmap(conf_matrix_best, annot=True, fmt='d', cmap='Blues', xticklabels=['No Use',
'Use'], yticklabels=['No Use', 'Use'])
plt.xlabel('Predicted')
plt.ylabel('Actual')
plt.title('Confusion Matrix - Best SVM')
plt.show()
```



k-Nearest Neighbors (kNN)

- Type: Instance-based Learning
- Best For:
 - Simple datasets with clear clusters or patterns.
 - When interpretability isn't a priority but simplicity is.
- Pros:
 - Easy to implement.
 - Makes no assumptions about the data distribution.
 - Adapts naturally to non-linear decision boundaries.
- Cons:
 - Computationally expensive for large datasets.
 - Sensitive to irrelevant features and feature scaling.
 - Performance depends heavily on the choice of k.

In [32]:

```
import numpy as np
from sklearn.neighbors import KNeighborsClassifier
from sklearn.metrics import classification_report, confusion_matrix, accuracy_score
import matplotlib.pyplot as plt
```

In [33]:

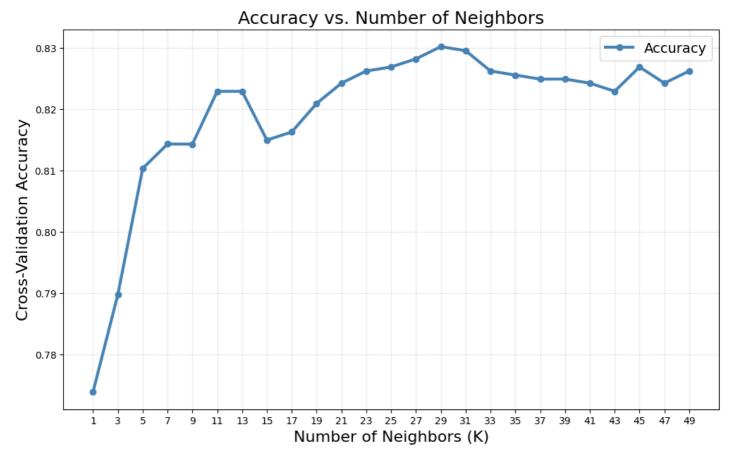
```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42
```

```
, stratify=y)
# Initialize KNN with K=3
knn = KNeighborsClassifier(n neighbors=3)
# Train the model
knn.fit(X train, y train)
# Predict on the test set
y pred = knn.predict(X test)
# Evaluate the model
print("Classification Report for K=3:")
print(classification report(y test, y pred))
# Confusion matrix
conf matrix = confusion matrix(y test, y pred)
print("Confusion Matrix for K=3:")
print(conf matrix)
Classification Report for K=3:
                          recall f1-score
                                              support
              precision
                   0.76
                             0.77
                                        0.76
                                                   156
                   0.84
                             0.83
                                       0.83
                                                   221
                                       0.80
                                                   377
    accuracy
   macro avg
                   0.80
                             0.80
                                       0.80
                                                   377
weighted avg
                   0.80
                             0.80
                                       0.80
                                                   377
Confusion Matrix for K=3:
[[120 36]
 [ 38 183]]
In [34]:
from sklearn.model selection import cross val score
# Test different values of K
acc = []
allks = range(1, 50, 2)
for k in allks:
   knn = KNeighborsClassifier(n neighbors=k)
    scores = cross val score(knn, X train, y train, cv=5, scoring='accuracy')
    acc.append(scores.mean())
    print(f"K = {k}, Cross-Validation Accuracy: {scores.mean():.3f}")
K = 1, Cross-Validation Accuracy: 0.774
K = 3, Cross-Validation Accuracy: 0.790
K = 5, Cross-Validation Accuracy: 0.810
K = 7, Cross-Validation Accuracy: 0.814
K = 9, Cross-Validation Accuracy: 0.814
K = 11, Cross-Validation Accuracy: 0.823
K = 13, Cross-Validation Accuracy: 0.823
K = 15, Cross-Validation Accuracy: 0.815
K = 17, Cross-Validation Accuracy: 0.816
K = 19, Cross-Validation Accuracy: 0.821
K = 21, Cross-Validation Accuracy: 0.824
K = 23, Cross-Validation Accuracy: 0.826
K = 25, Cross-Validation Accuracy: 0.827
K = 27, Cross-Validation Accuracy: 0.828
K = 29, Cross-Validation Accuracy: 0.830
K = 31, Cross-Validation Accuracy: 0.830
K = 33, Cross-Validation Accuracy: 0.826
K = 35, Cross-Validation Accuracy: 0.826
K = 37, Cross-Validation Accuracy: 0.825
K = 39, Cross-Validation Accuracy: 0.825
K = 41, Cross-Validation Accuracy: 0.824
K = 43, Cross-Validation Accuracy: 0.823
K = 45, Cross-Validation Accuracy: 0.827
K = 47, Cross-Validation Accuracy: 0.824
```

```
K = 49, Cross-Validation Accuracy: 0.826
```

In [35]:

```
# Plot accuracy vs K
plt.figure(figsize=(12, 7))
plt.plot(allks, acc, marker="o", color="steelblue", lw=3, label="Accuracy")
plt.xlabel("Number of Neighbors (K)", fontsize=16)
plt.ylabel("Cross-Validation Accuracy", fontsize=16)
plt.title("Accuracy vs. Number of Neighbors", fontsize=18)
plt.xticks(range(1, 50, 2))
plt.grid(alpha=0.25)
plt.legend(fontsize=14)
plt.show()
```



In [36]:

Classification Report for K=15:

0

precision

0.78

```
# Optimal K (max accuracy)
optimal_k = np.argmax(acc) + 1  # Adding 1 because index starts at 0
print(f"Optimal K: {optimal k}")
# Train KNN with the optimal K
knn optimal = KNeighborsClassifier(n neighbors=optimal k)
knn optimal.fit(X train, y train)
# Predict on the test set
y_pred_optimal = knn_optimal.predict(X test)
# Evaluate the optimal KNN model
print(f"Classification Report for K={optimal_k}:")
print(classification_report(y_test, y_pred_optimal))
# Confusion matrix for optimal K
conf matrix optimal = confusion matrix(y test, y pred optimal)
print(f"Confusion Matrix for K={optimal k}:")
print(conf_matrix optimal)
Optimal K: 15
```

recall f1-score

0.80

156

0.82

```
0.87
                              0.83
                                         0.85
                                                    221
                                         0.83
                                                     377
    accuracy
                    0.82
                              0.83
                                         0.82
                                                     377
   macro avg
                              0.83
                                         0.83
                                                    377
weighted avg
                    0.83
Confusion Matrix for K=15:
[[128 28]
 [ 37 184]]
```

In [37]:

```
import seaborn as sns

# Confusion matrix heatmap
sns.heatmap(conf_matrix_optimal, annot=True, fmt='d', cmap='Blues', xticklabels=['No Use', 'Use'], yticklabels=['No Use', 'Use'])
plt.xlabel('Predicted')
plt.ylabel('Actual')
plt.title(f'Confusion Matrix for K={optimal_k}')
plt.show()
```

