What is regression?

- Statistical models to explore the relationship between a response variable and some explanatory variables.
- Given values of explanatory variables, you can predict the values of the response variable.
- Response variable (a.k.a. dependent variable) The variable that you want to predict.
- Explanatory variables (a.k.a. independent variables) The variables that explain how the response variable will change.

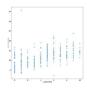
Visualizing two numeric variables

Before you can run any statistical models, it's usually a good idea to visualize your dataset. Here, you'll look at the relationship between house price per area and the number of nearby convenience stores using the Taiwan real estate dataset.

One challenge in this dataset is that the number of convenience stores contains integer data, causing points to overlap. To solve this, you will make the points transparent.

taiwan_real_estate is available from Taiwan_real_estate2.csv

- Import the seaborn and matplotlib packages
- Using taiwan_real_estate, draw a scatter plot of "price_twd_msq" (y-axis) versus "n_convenience" (x-axis).
- Draw a trend line calculated using linear regression. Omit the confidence interval ribbon. Note: The scatter_kws argument in regplot, scatter_kws={'alpha': 0.5}, makes the data points 50% transparent.





Transforming variables

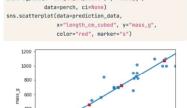
 Sometimes, the relationship between the explanatory variable and the response variable may not be a straight line. To fit a linear regression model, you may need to transform the explanatory variable or the response variable, or both of them.

Modeling mass vs. length cubed

```
perch["length_cm_cubed"] = perch["length_cm"] ** 3

mdl_perch = ols("mass_g ~ length_cm_cubed", data=perch).fit()
mdl_perch.params
```

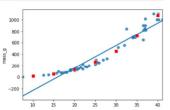




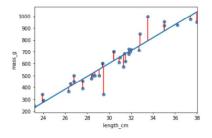
sns.regplot(x="length_cm_cubed", y="mass_g",

fig = plt.figure()





Residual standard error (RSE)



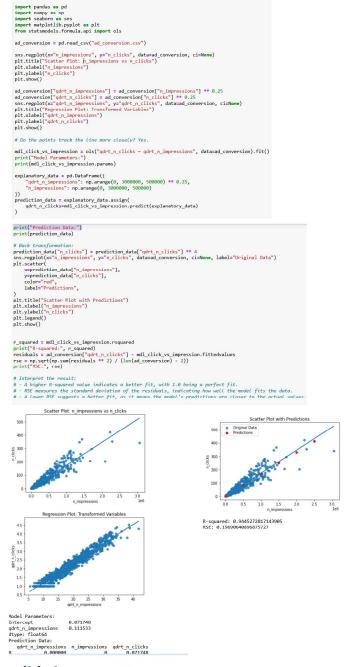
- A "typical" difference between a prediction and an observed response
- It has the same unit as the response variable.
- MSE = RSE²

Using scikit-learn to fit a classifier

```
from sklearn.neighbors import KNeighborsClassifier
X = churn_df[["total_day_charge", "total_eve_charge"]].values
y = churn_df["churn"].values
print(X.shape, y.shape)
```

(3333, 2), (3333,)

knn = KNeighborsClassifier(n_neighbors=15)
knn.fit(X, y)



scikit-learn syntax

```
from sklearn.module import Model
model = Model()
model.fit(X, y)
predictions = model.predict(X_new)
print(predictions)
```

array([0, 0, 0, 0, 1, 0])

Classifying labels of unseen data

- 1. Build a model
- 2. Model learns from the labeled data we pass to it
- 3. Pass unlabeled data to the model as input
- 4. Model predicts the labels of the unseen data
- Labeled data = training data

k-Nearest Neighbors

- Predict the label of a data point by
 - Looking at the k closest labeled data points
 - Taking a majority vote

```
#Task 1
import pandas as pd
import numpy as np
from sklearn.model selection import train test split
 from sklearn.neighbors import KNeighborsRegressor
from sklearn.metrics import mean_squared_error
from sklearn.model_selection import GridSearchCV
from sklearn.preprocessing import MinMaxScaler, LabelEncoder
# Load data
cc_apps = pd.read_csv("cc_approvals.data", header=None)
# Data Preprocessing
# Handling Missing Values (replace '?' with NaN)

cc_apps.replace('?', np.nan, inplace=True)

cc_apps.fillna(cc_apps.median(), inplace=True) # Filling missing values with median for num
# Convert non-numeric columns to strings
for column in cc_apps.columns:
    if cc_apps[column].dtype == 'object':
             cc_apps[column] = cc_apps[column].astype(str)
# Label Encoding for non-numeric columns
label_encoders = {}
for column in cc_apps.columns:
    if cc_apps[column].dtype == 'object':
        label_encoders[column] = LabelEncoder()
             cc_apps[column] = label_encoders[column].fit_transform(cc_apps[column])
cc_apps = cc_apps.drop([11, 13], axis=1)
# Split data into features and target variable
      cc_apps.drop(columns=[2]) # Replace 'target_column' with your actual target column name
y = cc_apps[2]
 # Split into train and test sets
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.33, random_state=42)
# Find the best k using Grid Search for KNeighborsRegressor
param_grid = {'n_neighbors': range(1, 21)} # Trying k values from 1 to 20
knn = KNeighborsRegressor()
grid_search = GridSearchCV(knn, param_grid, cv=5)
grid_search.fit(X_train_scaled, y_train)
# Print the best k value
print("Best K: ", grid_search.best_params_['n_neighbors'])
# Train the final KNN model with the best k value
best k = grid search.best params ['n_neighbors']
knn = KNeighborsRegressor(n_neighbors=best_k)
knn.fit(X_train_scaled, y_train)
 test_accuracy = knn.score(X_test, y_test)
print("Accuracy of the best KNN model: ", test_accuracy)
C:\Users\Lut Lat Aung\AppData\Local\Temp\ipykernel_29180\863855895.py:17: FutureWarning: Dro
ame reductions (with 'numeric_only=None') is deprecated; in a future version this will raise
mns before calling the reduction.
cc_apps.fillna(cc_apps.median(), inplace=True) # Filling missing values with median for n
Best K: 18
Accuracy of the best KNN model - -0 43739902524631247
```

```
import pandas as pd
cc_apps = pd.read_csv("cc_approvals.data", header = None)
print(cc_apps.head())
print(cc_apps.describe())
print('\n')
print(cc_apps.info())
print('\n')
cc_apps.tail(17)
cc_apps.fillna(cc_apps.mean(), inplace=True)
print(cc_apps.isnull().sum())
import numpy as np
cc_apps = cc_apps.replace('?', np.NaN)
for col in cc_apps.columns:
    if cc_apps[col].dtypes == 'object':
        cc_apps = cc_apps.fillna(cc_apps[col].value_counts().index[0])
print(cc_apps.isnull().sum())
print(cc_apps.describe())
print('\n')
print(cc_apps.info())
print('\n')
cc_apps.tail(17)
cc_apps = pd.get_dummies(cc_apps)
print(cc apps)
from sklearn.preprocessing import MinMaxScaler
X_train, y_train = cc_apps.iloc[:, :-1].values, cc_apps.iloc[:, [-1]].values
X_test, y_test = cc_apps.iloc[:, :-1].values, cc_apps.iloc[:, [-1]].values
scaler = MinMaxScaler(feature_range=(0, 1))
rescaledX_train = scaler.fit_transform(X_train)
rescaledX_test = scaler.transform(X_test)
from sklearn.linear_model import LogisticRegression
logreg = LogisticRegression()
logreg.fit(rescaledX_train,y_train)
```

Assessing classification performance

Predicted:	Predicted:
Legitimate	Fraudulent

Actual: Legitimate
Actual: Fraudulent

True Negative	
False Negative	True Positive

· Accuracy:

$$\frac{tp+tn}{tp+tn+fp+fn}$$

Precision

Predicted: Predicted: Fraudulent

Actual: Legitimate
Actual: Fraudulent

True Negative False Positive False Negative True Positive

Precision

$\frac{true\ positives}{true\ positives + false\ positives}$

Usually, the class of interest is called the positive class. As we aim to detect fraud, the positive class is an illegitimate transaction. So why is the confusion matrix important? There are other important metrics we can calculate from the confusion matrix. Precision is the number of true positives divided by the sum of all positive predictions, it is also called the positive predictive value. In our case, this is the number of correctly labeled fraudulent transactions divided by the total number of transactions classified as fraudulent. High precision means having a lower false positive rate.

Recall

Predicted: Predicted: Legitimate Fraudulent

Actual: Legitimate
Actual: Fraudulent

True Negative False Positive False Negative True Positive

Recall

$\frac{true\ positives}{true\ positives + false\ negatives}$

- High recall = lower false negative rate
- · High recall: Predicted most fraudulent transactions correctly

Recall is the number of true positives divided by the sum of true positives and false negatives. This is also called sensitivity. High recall reflects a lower false negative rate. For our classifier, it means predicting most fraudulent transactions correctly.

F1 score

• F1 Score: $2 * \frac{precision * recall}{precision + recall}$

The F1-score is the harmonic mean of precision and recall. This metric gives equal weight to precision and recall, therefore it factors in both the number of errors made by the model and the type of errors. The F1 score favors models with similar precision and recall, and is a useful metric if we are seeking a model which performs reasonably well acros both metrics.

```
from sklearn.model_selection import train_test_split

print(cc_apps.corr())
cc_apps_train, cc_apps_test = train_test_split(cc_apps, test_size=0.33, random_state=42)

from sklearn.metrics import confusion_matrix

y_pred = logreg.predict(rescaledX_test)
print("Accuracy of logistic regression classifier: ", logreg.score(rescaledX_test,y_test))
confusion_matrix(y_test,y_pred)

from sklearn.neighbors import KNeighborsClassifier
from sklearn.metrics import accuracy_score

for k in range(1, 10):
    knn_classifier = KNeighborsClassifier(n_neighbors=k)
    knn_classifier = KNeighborsClassifier(n_neighbors=k)
    knn_classifier.fit(X_train, y_train)
    y_pred = knn_classifier.predict(X_test)
    accuracy = accuracy_score(y_test, y_pred)
    print(f'k = {k}: Accuracy = {accuracy:.2f}')
```

k-Nearest Neighbors: Fit

In this exercise, you will build your first classification model using the churn_df dataset, can be loaded from churn df.csv.

The features to use will be "account_length" and "customer_service_calls". The target, "churn", needs to be a single column with the same number of observations as the feature data.

You will convert the features and the target variable into NumPy arrays, create an instance of a KNN classifier,

- Import KNeighborsClassifier from sklearn.neighbors.
- Create an array called X containing values from the "account_length" and "customer_service_calls" columns, and an array called y for the values of the "churn" column.
- Instantiate a KNeighborsClassifier called knn with 6 neighbors.
- Fit the classifier to the data using the .fit() method.

predict the labels of a set of new data points:

```
X = churn_df[['total_day_charge', 'total_eve_charge']].values
y = churn_df['churn'].values
print(X.shape, y.shape)
km = NNeighborsclassifier(n_neighbors=15)
[50.1, 10.9]])
print(X_new.shape)
predictions = knn.predict(X_new)
print(f'Predictions: {predictions}')
```

k-Nearest Neighbors: Predict

Now you have fit a KNN classifier, you can use it to predict the label of new data points. All available data was used for training, however, fortunately, there are new observations available, X_new.

The model knn, which you created and fit the data in the last exercise, will be used. You will use your classifier to

X new = np.array([[30.0, 17.5],[107.0, 24.1], [213.0, 10.9]])

```
X = churn_df[['total_day_charge', 'total_eve_charge']].values
y = churn_df['churn'].values
print(X.shape, y.shape)
knn = KNteighborsclassifier(n_neighbors=15)
```

Instructions

- · Create y_pred by predicting the target values of the unseen features
- Print the predicted labels for the set of predictions.

```
Predictions: [0 1 0]
```

```
import matplotlib.pyplot as plt
train_accuracies = {}
test_accuracies = {}
neighbors = np.arange(1,26)
print(neighbors)
for neighbor in neighbors:
         nt(neighbors)

neighbor in neighbors:

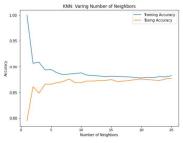
knn = KNeighborsClassifier(n_neighbors = neighbor)

knn.fit(X_train, y_train)

train_accuracies[neighbor] = knn.score(X_train, y_train)

test_accuracies[neighbor] = knn.score(X_test, y_test)
mprint(trut_accuractes.values())
print(test_accuracies.values())
my_train = list(train_accuracies.values())
my_test = list(test_accuracies.values())
plt.figure(figsize=(8,6))
plt.title("KNN: Varing Number of Neighbors")
plt.plot(neighbors, my_train, label="Training Accuracy")
plt.plot(neighbors, my_test, label="Tesing Accuracy")
plt.legend()
plt.xlabel("Number of Neighbors")
plt.ylabel("Accuracy")
plt.show()
```

[1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25] dict_values([0.795, 0.861, 0.864, 0.866, 0.866, 0.866, 0.871, 0.876, 0.869, 0.869, 0.872, 0.872, 0.873, 0.873, 0.875, 0.876, 0.877, 0.877, 0.873, 0.875, 0.875, 0.874, 0.873, 0.876, 0.877])



import pandas as pd sales df = pd.read csv('sales df.csv') X = sales_df['radio'].values
y = sales_df['sales'].values X = X.reshape(-1, 1)
print("Shape of X:", X.shape)
print("Shape of y:", y.shape) Shape of X: (4546, 1) Shape of y: (4546,)

from sklearn.linear_model import LinearRegression
reg = LinearRegression()
reg.fit(X_bmi, y)
predictions = reg.predict(X_bmi)
plt.scatter(X_bmi, y, color="Red")
plt.scatter(X_bmi, y, color="Red")
plt.ylabel("Blood Glucose (mg/dl)")
plt.xlabel("Blood Slucose (mg/dl)")
plt.xlabel("Body Mass Index")
nlt.show() plt.show()

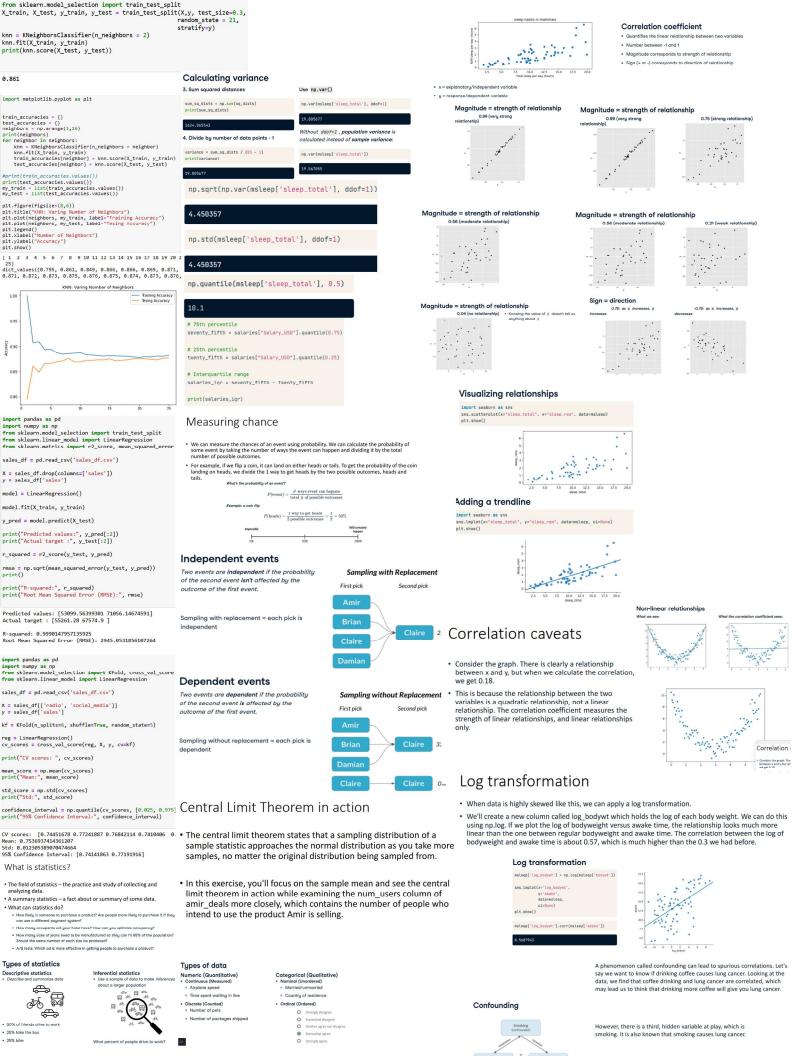
from sklearn.linear model import LinearRegression model = LinearRegression()
model.fit(X, y) predictions = model.predict(X) print("Five prediction values:", predictions[:5])

Five prediction values: [95491.17119147 117829.51

180 160 140 흥 120 100 60 400 Body Mass Index

111137.28167129] import matplotlib.pyplot as plt
plt.scatter(X, y, color='blue', label='Actual Observations') plt.plot(X, predictions, color='red', label='Linear Regression Model') plt.xlabel('Radio Advertising Expenditure')
plt.ylabel('Sales')
plt.legend(loc='best')
plt.show()

```
from sklearn.model_selection import train_test_split
 print(cc_apps.corr())
 #Drop the features 11 and 13
 cc_apps = cc_apps.drop([11, 13], axis=1)
 # Split into train and test sets
 cc_apps_train, cc_apps_test = train_test_split(cc_apps, test_size=0.33, random_state=42)
     1.000000 0.298902 0.271207 0.123121
     0.298902 1.000000 0.322330
                                       0.051345
 10
     0 271207
                0.322330
                           1 999999
                                       9 963692
     0.123121
                0.051345
                           0.063692
                                      1.000000
 # Import numpy
 import numpy as np
 # Replace the '?'s with NaN in the train and test sets
 cc_apps_train = cc_apps_train.replace('?', np.NaN)
cc_apps_test = cc_apps_test.replace('?', np.NaN)
 # Impute the missing values with mean imputation
 cc_apps_train.fillna(cc_apps_train.mean(), inplace=True)
 cc_apps_test.fillna(cc_apps_train.mean(), inplace=True)
 # Count the number of NaNs in the datasets and print the counts to verify
 print(cc_apps_train.isnull().sum())
 print(cc_apps_test.isnull().sum())
 for col in cc_apps_train.columns: # Iterate over each column of cc_apps_train
     if cc_apps_train[col].dtypes == 'object': # Check if the column is of object type
          # Impute with the most frequent value
# The value_counts() function returns a Series that contain counts of unique val
          # descending order so that its first element will be the most frequently-occurre
cc_apps_train = cc_apps_train.fillna(cc_apps_train[col].value_counts().index[0])
          cc_apps_test = cc_apps_test.fillna(cc_apps_train[col].value_counts().index[0])
# Count the number of NaNs in the dataset and print the counts to verify
print(cc_apps_train.isnull().sum())
print(cc_apps_test.isnull().sum())
 # At this point, there is no missing values.
# Convert the categorical features in the train and test sets independently
print(cc apps train)
cc_apps_train = pd.get_dummies(cc_apps_train)
cc_apps_test = pd.get_dummies(cc_apps_test)
print(cc_apps_train)
# Reindex the columns of the test set aligning with the train set
cc_apps_test = cc_apps_test.reindex(columns=cc_apps_train.columns, fill_value=0)
 # Import MinMaxScaler
from sklearn.preprocessing import MinMaxScaler
 # Searegate features and labels into separate variables
X_train, y_train = cc_apps_train.iloc[:, :-1].values, cc_apps_train.iloc[:, [-1]].values
X_test, y_test = cc_apps_test.iloc[:, :-1].values, cc_apps_test.iloc[:, [-1]].values
# Instantiate MinMaxScaler and use it to rescale X_train and X_test
scaler = MinMaxScaler(feature range=(0, 1))
 rescaledX_train = scaler.fit_transform(X_train)
rescaledX test = scaler.transform(X test)
# Import LogisticRegression
from sklearn.linear_model import LogisticRegression
 # Instantiate a LogisticRegression classifier with default parameter values
logreg = LogisticRegression()
# Fit logreg to the train set
logreg.fit(rescaledX_train,y_train)
# Import confusion_matrix
from sklearn.metrics import confusion_matrix
# Use logreg to predict instances from the test set and store it
y_pred = logreg.predict(rescaledX_test)
# Get the accuracy score of Logreg model and print it
print("Accuracy of logistic regression classifier: ", logreg.score(rescaledX_test,y_test))
# Print the confusion matrix of the logreg model
confusion_matrix(y_test,y_pred)
Accuracy of logistic regression classifier: 1.0
array([[103, 0],
[ 0, 125]], dtype=int64)
```



Nominal (Unordered) Married/unmarried (1/8)

Being able to identify data types is important since the type of data

Ordinal (Ordered)

Categorical data can be represented as numbers

- Somewhat agree (4) Strongly agree (5)

In reality, it turns out that coffee does not cause lung cancer and is only associated with it, but it appeared causal due to the third variable, smoking. This third variable is called a confounder, or lurking variable. This means that the relationship of interest between coffee and lung cancer is a spurious correlation.