

Logistic Regression



Agenda

- Types of Supervised Learning
- Standard process for Data Science Project
- Visiting Basics
 - Odds Ratio
- Binomial Logistic Regression
 - Maximum Likelihood Estimation
 - Assumptions of Logistic Regression
 - Significance of Coefficients



Agenda

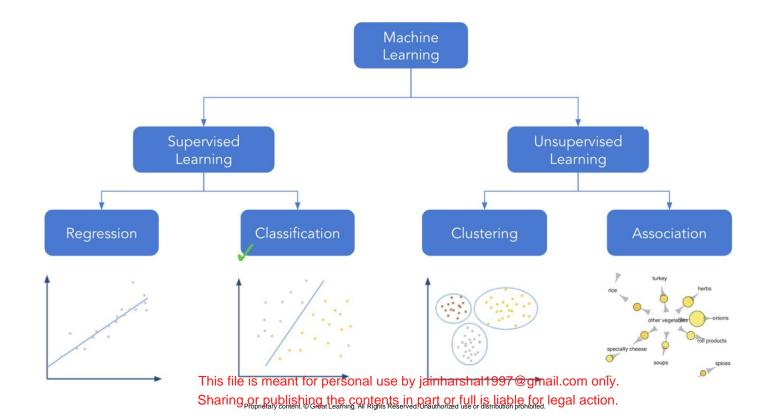
- Model Evaluation Metrics
 - Deviance
 - o AIC
 - Pseudo R²
- Model Performance Measures
 - Confusion Matrix
 - Cross Entropy
 - ROC AUC Score
- Imbalanced Data



Machine Learning



In this session, we shall cover classification



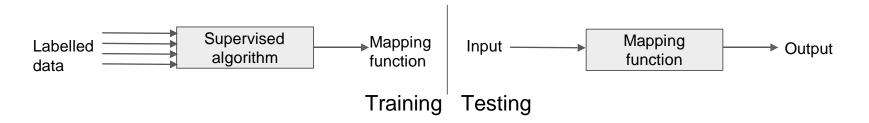


Supervised Learning



Supervised learning

Supervised learning aims at finding a model that maps the output (target) variable to the input (predictor) variables



Example: Detection of phishing emails based on certain phrases like 'You have won million'. More such phrases are prespecified while training the model. So if a new email also contains a similar phrase such emails can directly be tagged as spam.



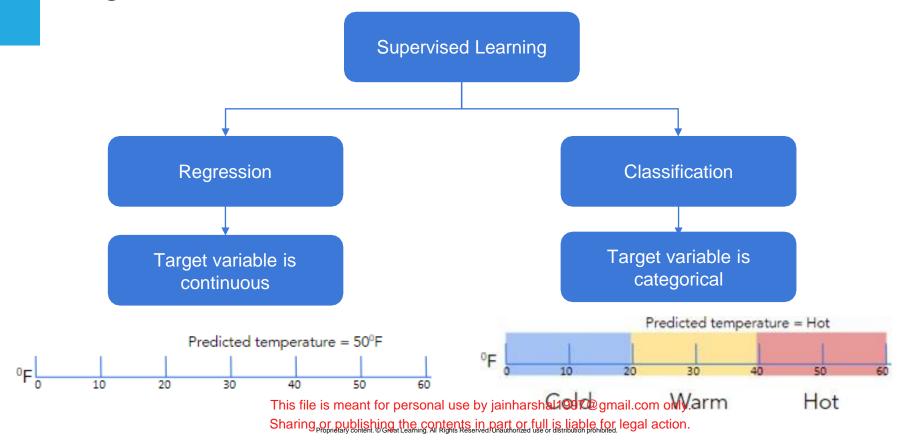
Supervised learning problems

Supervised learning is mainly used for two types of problems:

- Regression problem
- Classification problem



Regression vs classification





Classification

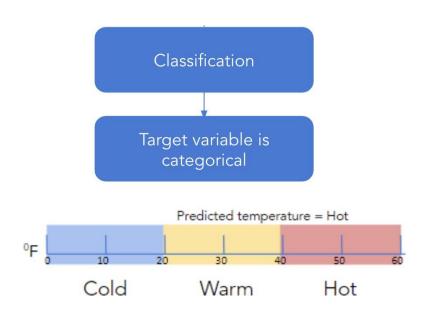


Class label

For classification, the target variable has categories.

In the example, Cold, Warm and Hot are the categories of the target variable.

These categories are called the class labels.





Classification

- An instance is mapped to one of many available labels
- Labels are the fixed number of values taken by the target variable
- The machine learns the pattern from train data where the labels are known for all instances. Then the learning can be used on new data where labels need to be predicted



Example of classification

Consider the example where we have inventory data for an online retailer which includes the number of orders, type and demand area for each item. The aim is to classify the items based on if they have a high or low demand.

This process can be automated using machine learning algorithms for classification.



Types of classification

Binary classification:

Classification with only two class labels

Example: Emails can be classified into spam and ham

Multiclass classification:

Classification with more than two distinct class labels

Example: Classification of land based on types of soils



Standard Process for Data Science Project



CRISP-DM

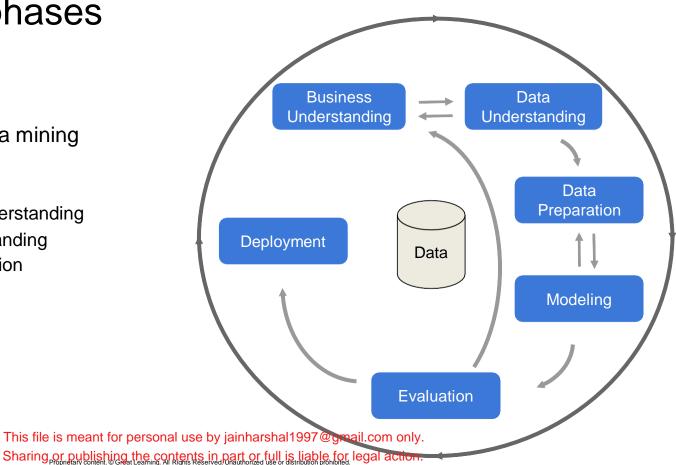
Cross Industry Standard Process for Data Mining (CRISP-DM) is a standard process used for data mining



CRISP-DM phases

CRISP-DM breaks data mining into six phases:

- **Business Understanding**
- **Data Understanding** 0
- **Data Preparation**
- Modeling
- **Evaluation**
- Deployment



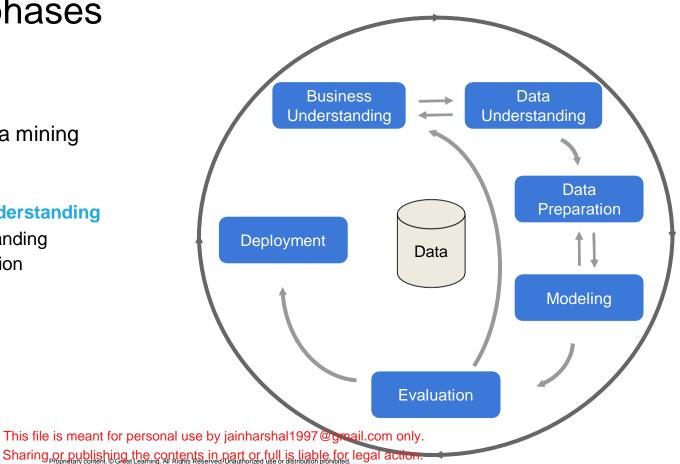
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CRISP-DM phases

CRISP-DM breaks data mining into six phases:

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- Deployment



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Business understanding

In this phase, we define what problem we are trying to solve.

Example: Consider the example where we have inventory data for an online retailer which includes the number of orders, type and demand area for each item. The aim is to classify the items based on if they have a high or low demand. So we can define clear business problems like:

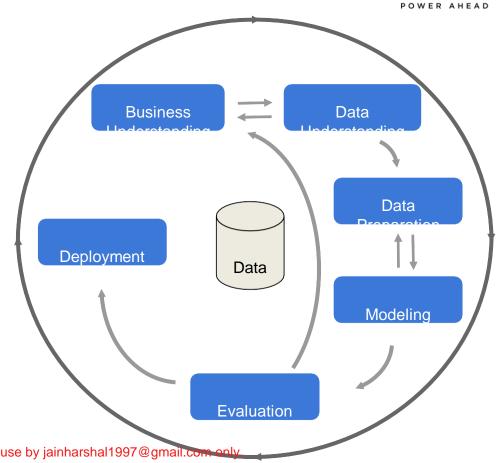
- Is the type of item related to the demand for the item?
- Can attributes in the considered data be used to classify the entire inventory list with reasonable accuracy?



CRISP-DM phases

CRISP-DM breaks data mining into six phases:

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Data understanding

This phase involves understanding the data considered for finding the solution.

Example: Consider the example where we have inventory data for an online retailer which includes the number of orders, type and demand area for each item. The aim is to classify the items based on if they have a high or low demand.

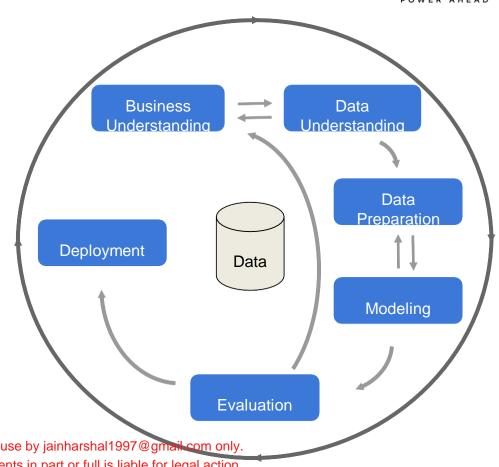
It is important to know if the items are perishable or non-perishable. For instance, items like dairy, cosmetics, and so on, can not be stocked, and hence adequate inventory should be available to meet the demand.



CRISP-DM phases

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Data preparation

This phase involves cleaning and processing the data to be in a format suitable for the model used to solve the problem.

Example: Consider the example to classify the items based on if they have a high or low demand. We can prepare the data as follows:

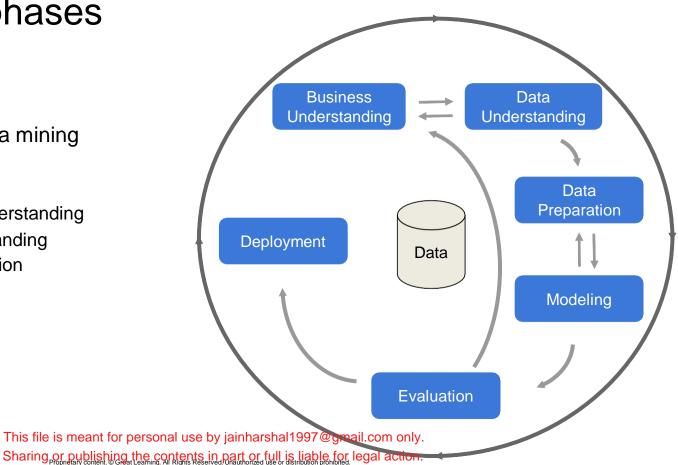
- Treat the missing values
- The categorical variables need to be dummy encoded
- Check for correlation among variables



CRISP-DM phases

CRISP-DM breaks data mining into six phases:

- **Business Understanding**
- **Data Understanding** 0
- **Data Preparation**
- **Modeling**
- **Evaluation**
- Deployment



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Modeling

 This phase involves finding the model that captures the solution to the business problem using available data

 We may have to try multiple models and go back and forth between data preparation and modelling to choose the correct model

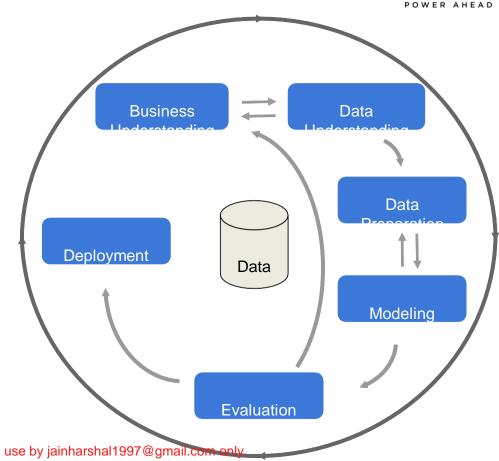
Example: In the modelling phase we try to find a function that maps the attributes like the number of orders, type, etc from the data to demand for the item.



CRISP-DM phases

CRISP-DM breaks data mining into six phases:

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Evaluation

Once the model is built we need to check how good the model performs on unseen data. This process is done during the evaluation phase.

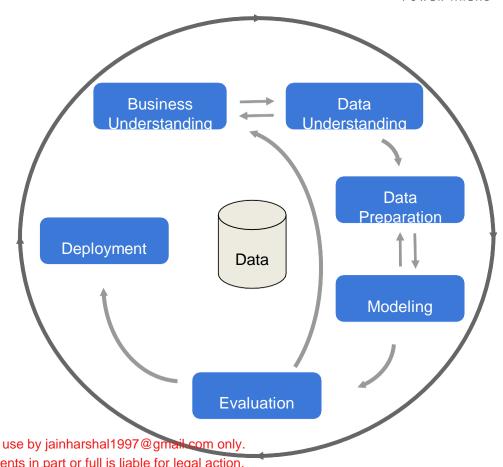
Example: We can check the model performance on data for which we know the actual demand. Using that data we can compare the predicted and actual values and evaluate.



CRISP-DM phases

CRISP-DM breaks data mining into six phases:

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Deployment

If we are satisfied with the performance of the model from the previous phase we deploy it in the deployment phase

Example: For the considered example of predicting demand for an item, perhaps we could develop an app that takes input as the attribute values for an item and returns the demand for that item to the retailer.



Visiting Basics



Odds vs probability

Odds of an event are the ratio of number of observations in favour of an event to number of observations not in favour of the event

Probability of an event is the ratio of number of observations in favour of an event to all possible observations

$$odds = \frac{\text{number of observations in favour of the event}}{\text{number of observations not in favour of the event}}$$

$$probability = \frac{\text{number of observations in favour of the event}}{\text{number of observations}}$$

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Odds vs probability

Plasma score	90	90	150	165	115	180	100	170	130	166
Is the patient Diabetic?	No	No	Yes	Yes	No	Yes	No	Yes	Yes	Yes

For the above data, the odds of a patient having diabetes is given by,

For the above data,, the probability of a patient having diabetes is given by,

odds =
$$\frac{\text{number of patients having diabetes}}{\text{number of patients not having diabetes}} = \frac{6}{4}$$
 probability = $\frac{\text{number of patients having diabetes}}{\text{Total number of patients}} = \frac{6}{10}$

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Log of odds

odds of having diabetes $=\frac{6}{4}$

log(odds of having diabetes) = ln (1.5) = 0.405

odds of not having diabetes $=\frac{4}{6}$

 $\log(\text{odds of not having diabetes}) = \ln(0.667) = -0.405$

Note: In=natural logarithm with base e.

- As we can see if we only consider odds value, the magnitude for each class value taken by variable is very different
- Hence, the log(odds) value is considered; so that no matter whichever the class is, the magnitude remains same
- Log of odds is the logit function used in logistic regression



Relation between odds and probability

If P(A) is probability of event A

$$Odds = \frac{P(A)}{1-P(A)}$$

$$Probability = \frac{odds}{1+odds}$$

$$log(Odds) = ln(\frac{P(A)}{1-P(A)})$$

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Odds ratio

- Odds ratio refers to the ratio of odds
- Odds ratio can be used to determine the impact of a feature on target variable
- For our considered example the odds ratio can be calculated as,

odds ratio =
$$\frac{\text{odds of patient having diabetes}}{\text{odds of patient not having diabetes}} = \frac{\frac{6}{4}}{\frac{4}{6}} = \frac{9}{4}$$





Question:

Patients with high sugar diet are considered more susceptible to diabetes. How can we determine whether sugar content in diet has an impact on possibility of a patient getting diagnosed with diabetes? Consider the following sample data.

Sugar content in diet	High	High	Low	High	Low	High	High	Low	High	Low
Is the patient Diabetic?	Yes	No	Yes	Yes	No	Yes	Yes	No	No	No





Solution:

From the given sample data we can calculate:

1. Odds of a patient having diabetes given he has high sugar diet

number of patients having diabetes given he has high sugar diet

number of patients not having diabetes given he has high sugar diet

2. Odds of a patient having diabetes given he has low sugar diet

number of patients having diabetes given he has low sugar diet

number of patients not having diabetes given he has low sugar diet

= 1

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Solution continued:

From 1 and 2 we can calculate odds ratio:

odds ratio =
$$\frac{\text{odds of a patient having diabetes given he has high sugar diet}}{\text{odds of a patient having diabetes given he has low sugar diet}} = \frac{\frac{4}{2}}{\frac{1}{3}} = 6$$

Thus from the odds ratio we can see that patients with a high sugar diet are 6 times more susceptible to diabetes compared to patients who have a low sugar diet.



Binomial Logistic Regression





Question:

Consider the example below about whether or not a patient has diabetes based on plasma score. Can we use linear regression line to predict the whether the patient is diabetic?

Plasma score	90	90	150	165	115	180	100	170	130	166
Is the patient Diabetic?	No	No	Yes	Yes	No	Yes	No	Yes	Yes	Yes





Answer:

The example about whether or not a patient has diabetes based on plasma score is a classification problem. Moreover, the target variable is categorical. Hence, we can not use linear regression line to predict the whether the patient is diabetic. We classify them as diabetic and non-diabetic.

Plasma score	90	90	150	165	115	180	100	170	130	166
Is the patient Diabetic?	No	No	Yes	Yes	No	Yes	No	Yes	Yes	Yes



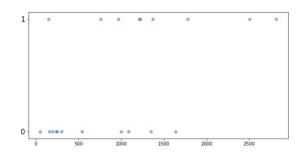
- Here on we shall consider the adjacent data
- The data tells us the presence of one fish depending on the density of other fish in a lake
- If they compete with each other, then the higher density of one may suggest the absence of the other whereas if they are symbiotic, high density of one may promote the other

BKT kg/ha	Presence of fish
1085.33	0
1210	1
1780.62	1
52.4	0
200	0
2502.67	1
301.33	0
542	0
969.33	1
240.56	0
1640	0
247	0
999.99	0
1220.76	1
150.67	1
160	0
2816	1
760	1
1350	0
1370	m only

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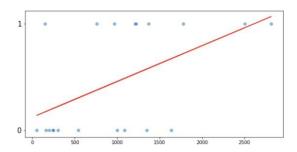


 Consider the scatter plot of the previous data (data on slide 42)



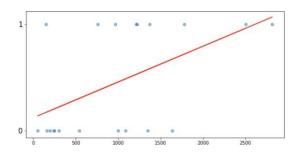
Fit a linear regression line to it

Note the line is not a true representative of the data

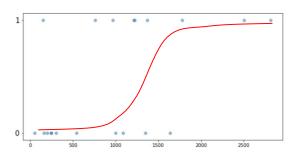




 A S-shaped curve as in the figure below gives the true relationship



Such a curve is given by the sigmoid function



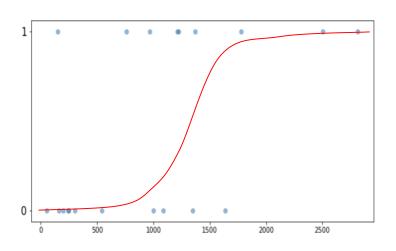


What is a sigmoid function?

 The sigmoid function is a mathematical function which is S-shaped and is given by

$$y(x)=rac{1}{1+e^{-(mx+b)}}$$

It exists between 0 to 1





- Logistic regression is a binary classification algorithm
- It predicts the probability of occurrence of a class label. Based on these probabilities the data points are labelled
- A threshold (or cut-off; commonly a threshold of 0.5 is used) is fixed, then

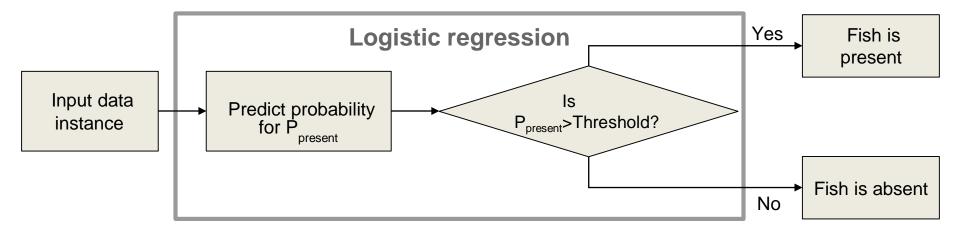
	Classify as
threshold < probability	Presence of fish
Probability > threshold	Absence of fish

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Main steps in logistic regression

Consider that logistic regression is used to identify whether or not a patient is suffering from diabetes



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The logistics regression is given by

$$\pi(x)=rac{e^{eta_0+\sum_{i=1}^neta_ix_i}}{1+e^{eta_0+\sum_{i=1}^neta_ix_i}}$$

- Here, π(x) is the conditional expectation of the outcome given the values for independent variables, i.e. E(Y|X)
- It predicts the probability of occurrence of a class label by fitting the data to a function called logit function, hence called logit regression



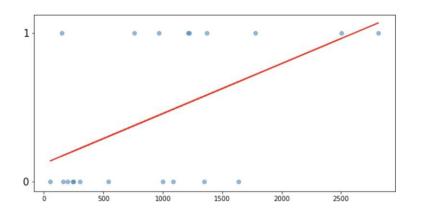
Probability as output of logistic regression

- The logistic regression model is given by $\pi(x)=rac{e^{eta_0+\sum_{i=1}^neta_ix_i}}{1+e^{eta_0+\sum_{i=1}^neta_ix_i}}$
- ullet Taking limits tending to - ∞ on both side, $\lim_{x o-\infty}rac{e^x}{1+e^x}=0$

ullet Taking limits tending to ∞ on both side, $\lim\limits_{x o\infty}rac{e^x}{1+e^x}=1$

• Thus $\pi(x)$ lies in-between 0 and 1, i.e. $\pi(x) \in [0,1]$ and can be viewed as probability





- Since we are predicting probabilities values need to be between 0 and 1
- Consider the points A and B for which values for plasma score are 80 and 177 respectively, probability values are out of the expected range 0 to 1
- Hence linear regression cannot be directly used to predict probabilities



Usage

Classification:

The ICU in a hospital is assigned on priority to high risk patients. Logistic regression can be used to classify the list of patients into high risk and low risk records.

Profiling:

Nuclear fuel companies moderate various factors like pressure, temperature, etc. to produce high yielding and low yielding fuels. Based on different parameters for the current day, it can be predetermined whether the fuel produced will be high yielding or not. The company can then alter the parameters to produce high yielding fuel every day.

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 Thus for a binary logistic classification where the target variable takes two values names 0 and 1, we have

Class labels	0	1
P[Y=y X]	1-π(x)	π(x)

- π(x) denotes the probability that the response is present for the records for some combination of values that the independent variables take, i.e. for X=x
- 1-π(x) denotes the probability that the response is absent for the records for some combination of values that the independent variables take i.e. for X=x



Linearization

 To estimate the parameter we need to linearize the function. We use the logit transformation

$$\eta=\lnrac{\pi}{1-\pi} \hspace{1cm}\pi(x)=rac{e^{eta_0+\sum_{i=1}^neta_ix_i}}{1+e^{eta_0+\sum_{i=1}^neta_ix_i}}$$

- The ratio $\pi/(1-\pi)$ is called the odds
- Hence the logit transformation is also known as the log-odds



Linearization

We have,

$$\frac{\pi(x)}{1-\pi(x)} = \frac{\frac{\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}{1+\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}}{1-\frac{\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}{1+\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}} = \frac{\frac{\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}{1+\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}}{\frac{1}{1+\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}} = \frac{\exp^{\beta_0 + \sum_{i=1}^n \beta_i x_i}}{1}$$

That is,

$$rac{\pi(x)}{1-\pi(x)} = \exp^{eta_0 + \sum_{i=1}^n eta_i x_i}$$
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Linearization

Taking natural log on both sides we have,

$$\ln rac{\pi(x)}{1-\pi(x)} = \ln \exp^{eta_0 + \sum_{i=1}^n eta_i x_i}$$

$$\ln rac{\pi(x)}{1-\pi(x)} = eta_0 + \sum_{i=1}^n eta_i x_i$$

Thus, we have a linear relationship.



Interpreting the parameter

The logistic regression is given by

$$\ln rac{\pi(x)}{1-\pi(x)} = eta_0 + \sum_{i=1}^n eta_i x_i$$

- In a linear regression model, β_1 gives the average change in Y associated with a one-unit increase in X
- Whereas, in a logistic regression model, increasing X by one unit changes the log odds by β₁, or equivalently it multiplies the odds by exp{β₁}



Interpreting for our example

The logistic regression is given by

$$\ln rac{\pi(x)}{1-\pi(x)} = 0.142 + 1.0018\,$$
 BKT kg/ha

Increasing BKT kg/ha by one unit changes the log odds by 1.0018

OR

Increasing BKT kg/ha by one unit multiplies the odds by exp{1.0018}
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BKT kg/ha	Presence of fish
1085.33	0
1210	1
1780.62	1
52.4	0
200	0
2502.67	1
301.33	0
542	0
969.33	1
240.56	0
1640	0
247	0
999.99	0
1220.76	1
150.67	1
160	0
2816	1
760	1
1350	0
1370	1



Parameter estimation

The general form of logistic regression model is

$$y_i = E(y_i) + \epsilon_i$$

where the observations yi are independent bernoulli random variables

The expected value of y_i's is

$$E(y_i) = \pi_i = rac{e^{eta_0 + \sum_{i=1}^n eta_i x_i}}{1 + e^{eta_0 + \sum_{i=1}^n eta_i x_i}}$$

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Parameter estimation

- The method of least squares fails
- Since, unlike the linear regression, the closed form solution does not exist

where a closed form solution is an exact solution evaluated with a fixed number of operations

So we use the method of maximum likelihood estimation





Question:

Consider the example below about whether or not a patient has diabetes based on plasma score. Since the logistic regression equation for this example will represent a straight line can we determine the residual values?

Plasma score	90	90	150	165	115	180	100	170
Is the patient Diabetic?	No	No	Yes	Ye s	No	Yes	No	Yes





Solution:

For the considered example the equation of logistic regression line is given by,

$$\ln\left(\frac{p}{1-p}\right) = \beta_0 + \beta_1$$
 Plasma Score

p = probability of a patient having diabetes

$$\ln\left(\frac{p}{1-p}\right)$$
 = logit function

 β_0 = Intercept value

 β_1 = Coefficient for variable Plasma Score

Plasma Score = Values, taken by variable Plasma Score @gmail.com only.





Solution continued:

The value of logit function i.e the LHS for a patient who has diabetes is given by

$$\ln\left(rac{p}{1-p}
ight) = \ln\left(rac{1}{0}
ight) = \ln(1) - \ln(0) = 0 - (-\infty) = +\infty$$
 ...(i)

The value of logit function i.e the LHS for a patient who has diabetes is given by

$$\ln\left(rac{p}{1-p}
ight) = \ln\left(rac{0}{1}
ight) = \ln(0) - \ln(1) = -\infty - 0 = -\infty$$
 ...(ii)





Solution continued:

From i and ii we can infer that the logistic regression line will approach $+\infty$ and $-\infty$ values. So the residuals will also take infinite values and hence cannot be determined.



Why we cannot use least squares for logistic regression?

- The residual values for the logistic regression line cannot be determined
- Least squares optimization cannot be performed without determining residuals for the line
- So we cannot use least squares optimization for logistic regression and instead use an optimization approach called maximum likelihood estimation
- Note: least squares for linear regression is a special case of maximum likelihood estimation



Maximum likelihood estimation

- The method of Maximum Likelihood Estimation (MLE) is a method of estimating the parameter of a function by maximizing the likelihood function
- The likelihood function is the joint probability density function of the sample
- For binomial logistic regression the data follows the bernoulli distribution. So the probability distribution is given by

$$f(y_i) = \pi_i^{y_i} (1 - \pi_i)^{1 - y_i}$$



Maximum likelihood estimation

The likelihood function (L) is given as

$$L = \prod_{i=1}^n f(y_i) = \prod_{i=1}^n \pi_i^{y_i} (1-\pi_i)^{1-y_i}$$

since we have an independent sample, L is the product of all probabilities

• Taking log on both sides, we have

$$\ln L = \ln \prod_{i=1}^n \pi_i^{y_i} (1 - \pi_i)^{1 - y_i}$$

$$\ln L = \sum_{i=1}^n y_i \ln \pi_i + \sum_{i=1}^n (1-y_i) \ln (1-\pi_i)$$
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Maximum likelihood estimation

Taking log on both sides, we have

$$\ln L = \sum_{i=1}^n y_i \pi_i + \sum_{i=1}^n (1-y_i)(1-\pi_i)$$

So solving, we have

$$\ln L = \sum_{i=1}^n [y_i \ln rac{\pi_i}{1-\pi_1}] + \sum_{i=1}^n \ln (1-\pi_i)$$

 This equation is further solved by numerical method such as the Newton-Raphson method in order to get the estimates



Assumptions of Logistic Regression



Assumptions

We have seen that the logistic regression can be linearized, so it has assumptions almost same as that of linear regression

Assumption 1	Independence of error, whereby all sample group outcomes are separate from each other (i.e., there are no duplicate responses)
Assumption 2	Linearity in the logit for any continuous independent variables
Assumption 3	Absence of multicollinearity
Assumption 4	lack of strongly influential outliers

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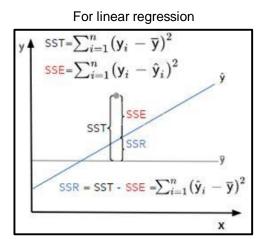


Significance of Coefficients



Significance of coefficients

- In a linear regression model, the significance of a regression coefficient is determined with the help of a t-test
- In logistic regression, the significance of the coefficients is determined by the wald statistic and by the likelihood ratio test
- To test the significance of the model, the likelihood ratio test is used





Significance of coefficients - Wald test

For β to be significant, β > 0.

against

$$H_0: \beta = 0$$
 against $H_1: \beta \neq 0$

It implies

$$H_0$$
: The parameter β is not significant H_1 : The parameter β is significant

• Failing to reject H₀ implies that the parameter β is not significant



Significance of coefficients - Wald test

The Wald statistic is given by

$$Z_{wald} = rac{\hat{eta}}{SE(\hat{eta})}$$
 where \hat{eta} is the estimated value of eta .

The Wald statistic follows the N(0,1) distribution

• Decision Rule: Reject H_0 if $|Z| > Z_{\alpha/2}$ or if the p-value is less than the α (level of significance)



Significance of coefficients - LRT

• For β to be significant, $\beta > 0$

$$H_0: \beta = 0$$
 against $H_1: \beta \neq 0$

It implies

 H_0 : The parameter β is not significant

against H_1 : The parameter β is significant

• Failing to reject H₀ implies that the parameter β is not significant



Significance of coefficients - LRT

The likelihood ratio test (LRT) is given by

$$D = -2 \ln \left[\frac{\text{Likelihood of model without predictors}}{\text{Likelihood of model with predictors}} \right]$$

- D follows the chi-squared distribution with one degree of freedom, i.e. χ₁
- Decision Rule: Reject H_0 if $\chi \ge \chi_{\alpha/2}$ or if the p-value is less than the α (level of significance)



Significance of model - LRT

 The hypothesis for testing all coefficient in logistic regression can be extended from the previous test for one coefficient as follows

$$H_0$$
: $\beta_1 = \beta_2 = \beta_3 = 0$ against H_1 : At least one $\beta_k \neq 0$ $(k = 1, 2, 3)$

It implies

H₀: the logistic model is not significant

against H₁: the logistic model is significant



Significance of model - LRT

The likelihood ratio test is given by

$$D = -2 \ln \left[\frac{\text{Likelihood of model without predictors}}{\text{Likelihood of model with predictors}} \right]$$

- D follows the chi-squared distribution with one degree of freedom, i.e. χ₁
- Decision Rule: Reject H_0 if $\chi \ge \chi_{\alpha/2}$ or if the p-value is less than the α (level of significance)



Model Evaluation Metrics



Model evaluation metrics

The model evaluation metrics are

- Deviance
- AIC
- Pseudo R²



Terminologies

- Null model: A model without any predictors
- Saturated model: A model with exactly n samples (n predictors), that fits the data perfectly
- Full model: A model fitted with all the variables in the data
- Fitted model: A model with at least one predictor variable



- Deviance is analogous to the sum of squares in the linear regression
- A measure of goodness of fit for a logistic regression
- Given by

$$D = -2 \ln \left[\frac{\text{Likelihood of fitted model}}{\text{Likelihood of saturated model}} \right]$$

where saturated model is a model assumed to have the perfect fit

If the saturated model is not available use the fitted model.



- Null deviance: The difference between the log likelihood of the null model and saturated model
- Model deviance: The difference between the log likelihood of the null model and fitted model
- Smaller values indicate a better fit
- To check for significance of k predictors, subtract the model deviance from the null deviance and access it on X_k



• In linear regression, we have R² defined as the ratio of explained variation to the total variation

$$R^2 = \frac{\text{Explained variation}}{\text{Total variation}} = \frac{\text{SSR}}{\text{SST}}$$

- In logistic regression, we can consider the deviance similar to the R² in linear regression
- Deviance can be thought of as the R² value such that the denominator is the total variation and the numerator is the variation explained by the fitted model

$$D = -2 \ln \left[rac{ ext{Likelihood of fitted model}}{ ext{Likelihood of saturated model}}
ight]$$



AIC

- The Akaike Information Criteria (AIC) is a relative measure of model evaluation for a given dataset
- It is given by:

$$AIC = -2 \ln L + 2K$$
 L: log-likelihood K: parameters to be estimated

 The AIC gives a trade-off between the model accuracy and model complexity, i.e. it prevents us from overfitting



Pseudo R²

- The non-pseudo R² or the R² in the linear regression framework is the explained variability and the correlation (for simple linear regression)
- An equivalent R² statistic does not exist in the logistic regression since the parameters are estimated by the method of maximum likelihood
- However, there are various pseudo R²s developed which are similar on the scale, i.e.
 on [0,1], and work exactly the same with higher values indicating a better fit



Pseudo R²

The pseudo R² are

McFadden R²

• Cox-Snell R²

Nagelkerke R²



McFadden R²

It is defined as

$$R_{ ext{McFadden}}^2 = 1 - rac{\ln ext{likelihood of full model}}{\ln ext{likelihood of null model}}$$

- If comparing two models on the same data, the model which has higher value is considered to be better
- The pseudo R2 in the python output is the McFadden R²



Cox-Snell R²

• It is similar to the McFadden R² and is defined as

$$R_{ ext{Cox-Snell}}^2 = 1 - \left\{ rac{ ext{ln likelihood of null model}}{ ext{ln likelihood of full model}}
ight\}^{rac{2}{N}}$$

- The likelihood is the product of probability N observations of the dataset. Thus the Nth square root of the provides an estimated of each target value
- The R² is less than 1
- For a model with likelihood 1, i.e the predictions are perfect, then the denominator becomes 1

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Nagelkerke R²

It is based on Cox-Snell R², it scales the values so that the maximum is 1

$$R_{ ext{Nagelkerke}}^2 = rac{1 - \left\{rac{\ln ext{likelihood of null model}}{\ln ext{likelihood of full model}}
ight\}^{rac{2}{N}}}{1 - \left\{\ln ext{likelihood of null model}
ight\}^{rac{2}{N}}}$$

- If the full model predicts the outcome perfectly, i.e it has likelihood = 1, then R²_{Nagelkerke} = 1
- Similarly, if likelihood of null model is equal to that of full model then R² = 0



Model Performance Measures



Performance metrics

The following metrics can be used to evaluate the performance of classification models:

- Confusion matrix
- Cross entropy
- ROC



Performance metrics

We shall consider the fish data as described before

BKT kg/ha	Presence of fish
1085.33	0
1210	1
1780.62	1
52.4	0
200	0
2502.67	1
301.33	0
542	0
969.33	1
240.56	0
1640	0
247	0
999.99	0
1220.76	1
150.67	1
160	0
2816	1
760	1
1350	0
1370	m only

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Performance metrics

The following metrics can be used to evaluate the performance of classification models:

- Confusion matrix
- Cross entropy
- ROC



Confusion matrix

Performance measure for classification problem

It is a table used to compare predicted and actual values of the target variable

		← —Actual	values		
		Positive(1)	Negative(0)		
d values 🕶	Positive(1)	True Positive: Predicted value is positive and the actual value is also positive	False Positive: Predicted value is positive but the actual value is negative		
	Negative(0)	False Negative: Predicted value is negative but the actual value is positive	True Negative: Predicted value is negative and the actual value is also negative		
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Confusion matrix for our example

Confusion matrix for our considered example to predict presence of fish is given as:

Predicted values

Positive(1) Positive(1) True Positive False Positive 3 Positive(1) False Negative Negative(0)

Actual values



Performance evaluation metrics

Confusion matrix can be used to calculate the following evaluation metrics for a model:

- Accuracy
- Precision
- Recall
- False Positive Rate
- Specificity
- F₁ score
- Kappa



Accuracy

Actual values

	Positive(1)	Negative(0)
Positive(1)	True Positive(TP)	False Positive(FP)
Negative(0)	False Negative(FN)	True Negative(TN)

 Accuracy is the fraction of predictions that out model got correct

$$Accuracy = \frac{\text{number of correctly predicted records}}{\text{Total number of records}}$$

Accuracy =
$$\frac{TP+TN}{TP+TN+FP+FN}$$

Higher the value of accuracy better is the model





Accuracy is not always a reliable metric

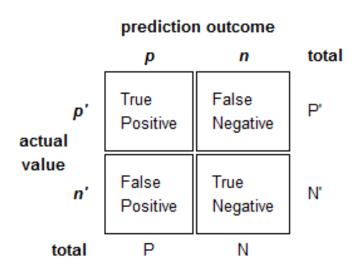
Consider a dataset with information about 1000 patients. 960 of those patients have diabetes and only 40 do not have diabetes. We have a model 'A' that classifies every patient as diabetic.

Accuracy of model A
$$=$$
 $\frac{960}{1000} = 96 \%$

Even though the accuracy for model A is high it is not a good model. Since even when it will encounter information about a new patient it will always predict that the patient is diabetic. This scenario when accuracy is not a reliable metric is called the the metric is called the the contents in part or full is liable for legal action.



Precision



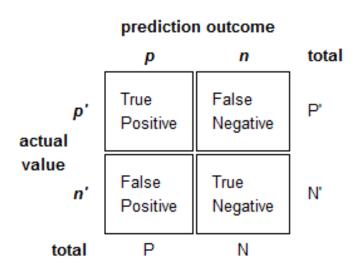
 Precision is proportion of positive cases that were correctly predicted

Precision =
$$\frac{TP}{TP+FP}$$

Higher is the precision better the model



Recall



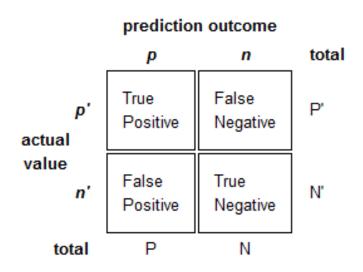
- Recall is the proportion of actual positive cases that were correctly predicted
- Recall is also sometimes called True Positive Rate (TPR) or Sensitivity

Recall =
$$\frac{TP}{TP+FN}$$

Higher value of TPR implies a better model



False positive rate



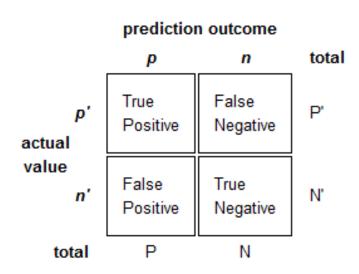
 False Positive Rate (FPR) is the proportion of actual negative cases that were predicted positive (incorrectly)

$$FPR = \frac{FP}{FP+TN}$$
 $FPR = 1 - Specificity$

Lower the value of FPR better is the model



Specificity



 Specificity is the proportion of actual negative cases that were correctly predicted

Specificity =
$$\frac{TN}{TN+FP}$$

Higher the specificity better is the model



F₁ score

- F₁ score is the harmonic mean of precision and recall values for a classification model
- It is a good measure if we want to find a balance between precision and recall or if there is an uneven distribution of classes (either positive or negative class has way more actual instances than the other)

$$F_1 ext{score} = \left(rac{ ext{recall}^{-1} + ext{precision}^{-1}}{2}
ight) = 2 \cdot rac{ ext{precision} \cdot ext{recall}}{ ext{precision} + ext{recall}}$$

Higher the F₁ score better the model



Performance metrics for our example

Performand Metric	e Accuracy	Precision	Recall	Specificity	False positive rate	F ₁ score
Formulae	$\frac{TP+TN}{TP+TN+FP+FN}$	$rac{TP}{TP+FP}$	$rac{TP}{TP+FN}$	$rac{TN}{TN+FP}$	1 – Specificity	$2 \cdot \frac{\text{precision} \cdot \text{recall}}{\text{precision} + \text{recall}}$
Value for o	0.7	0.667	0.667	0.727	0.273	0.667



Reliability

- Reliability is the degree to which an assessment tool produces consistent results
- Inter-rater reliability is used to measure the degree to which different raters agree while assessing the same thing
- In the case of logistic regression, the raters are the actual labels and predicted labels for the categorical target variable
- In logistic regression, the inter-rater reliability is the number of labels that match in both the predicted and actual instances



Reliability - Kappa statistic

The kappa statistics is used to test inter-rater reliability

$$\kappa = rac{p_o - p_e}{1 - p_e}$$

 p_o = relative observed agreement between raters

p_e = hypothetical probability of chance agreement



Kappa statistic

 Kappa statistic is a measure of interrater reliability or degree of agreement

 Kappa statistic can take values from the range [-1,1]

Карра	Interpretation
<0	No agreement
0-0.2	Slight agreement
0.2-0.4	Fair agreement
0.4-0.6	Moderate agreement
0.6-0.8	Substantial agreement
0.8-1	Almost perfect agreement

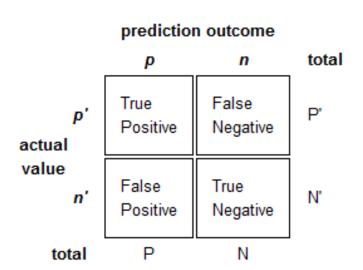


Calculation of kappa

- Calculate p_o, let A: Actual values and B: Predicted values
- Calculate P(A ∩ B)_{positive}
 - P(A ∩ B)_{positive} = Probability that both actual and predicted values are positive
- Calculate P(A ∩ B)_{negative}
 - P(A ∩ B)_{negative} = Probability that both actual and predicted values are negative
- 1. $p_e = P(A \cap B)_{positive} + P(A \cap B)_{negative}$



Calculation of p_o



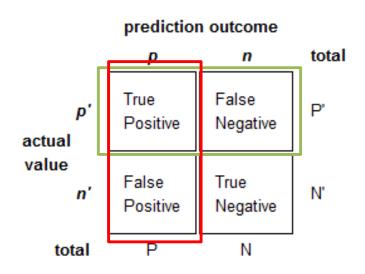
p_o is the observed agreement i.e when the actual and predicted values match

$$p_o = rac{ ext{number of instances in agreement}}{ ext{total instances}}$$

$$p_o = rac{TP + TN}{TP + TN + FP + FN}$$



Calculation of $P(A \cap B)_{positive}$



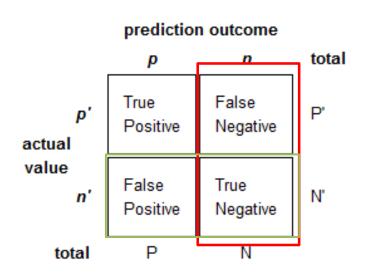
Since A and B are independent events:

$$P(A \cap B)_{positive} = P(A)_{positive} P(B)_{positive}$$

P(A) _{positive}	P(B) _{positive}
$\frac{TP{+}FN}{TP{+}TN{+}FP{+}FN}$	$\frac{TP{+}FP}{TP{+}TN{+}FP{+}FN}$



Calculation of $P(A \cap B)_{negative}$



Since A and B are independent events:

$$P(A \cap B)_{\text{negative}} = P(A)_{\text{negative}}$$
. $P(B)_{\text{negative}}$

P(A) _{negative}	P(B) _{negative}
$\frac{FP + TN}{TP + TN + FP + FN}$	$rac{FN + TN}{TP + TN + FP + FN}$



Calculation of p_e

p_e is hypothetical probability of chance agreement i.e when either both actual and predicted values are positive or both are negative

$$p_e = P(A \cap B)_{\text{positive}} + P(A \cap B)_{\text{negative}}$$



Performance metrics

The following metrics can be used to evaluate the performance of classification models:

- Confusion matrix
- Cross entropy
- ROC



Cross Entropy

- Cross entropy is the loss function commonly used in classification problems
- As the prediction goes closer to actual value the cross entropy decreases

$$H(y) = -\sum_{i} \mathsf{y}_{act(i)} \ln \left(\mathsf{y}_{pred(i)}
ight)$$

i = class (0 or 1)

H(y) = cross entropy

 $y_{act(i)}$ = actual probability for class i

 $y_{\text{pred(i)}} = \underset{\text{This file is meant for personal use by jainharshal1997@gmail.com only.}}{\text{This file is meant for personal use by jainharshal1997@gmail.com only.}}$



Calculation of cross entropy

Let us consider one observation from our considered example:

	Actual probability	Predicted probability
Fish present	1	0.72
Fish not present	0	0.28

$$egin{aligned} H(y) &= &- \sum_{i} \mathsf{y}_{act(i)} \ln \left(\mathsf{y}_{pred(i)}
ight) \ &= &- \mathsf{y}_{act(1)} \ln \left(\mathsf{y}_{pred(1)}
ight) - \mathsf{y}_{act(0)} \ln \left(\mathsf{y}_{pred(0)}
ight) \ &= &- 1 \cdot \ln (0.72) - 0 \cdot \ln (0.28) \ &= &- (\, -0.3285) = 0.33 \end{aligned}$$



Performance metrics

The following metrics can be used to evaluate the performance of classification models:

- Confusion matrix
- Cross entropy
- ROC

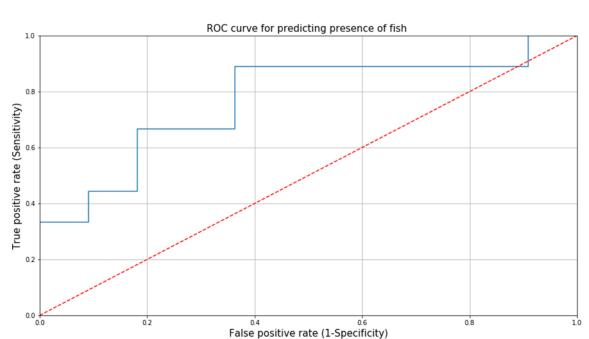


ROC

- Receiver operating characteristics (ROC) curve is
- The TPR and FPR values change with different threshold values
- ROC curve is the plot of TPR against the FPR values obtained at all possible threshold values



ROC curve for our example



TPR	FPR	Threshold
0.33	0	0.78
0.33	0.09	0.73
0.44	0.091	0.62
0.44	0.18	0.61
0.66	0.18	0.55
0.67	0.36	0.46
0.88	0.36	0.35
0.89	0.91	0.159
1	0.91	0.157

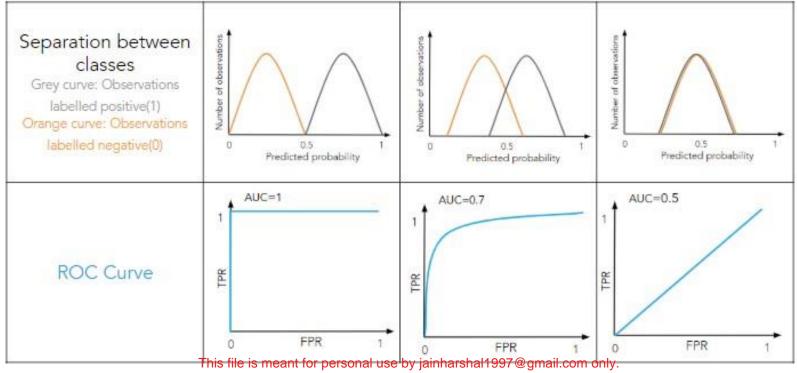


AUC

- Area under the ROC curve (AUC) is the measure of separability between the classes of target variables
- AUC increases as the separation between the classes increases
- Higher the AUC better the model



Effect of separation between classes on ROC



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Youden's index

- Sensitivity and specificity represent the total number of correctly identifies samples (true positives and the true negatives)
- Youden's index is the classification cut-off probability for which the (Sensitivity + Specificity -1) value is maximized
- Higher the value of Youden's index better the model

Youden's Index =
$$\max$$
 (Sensitivity + Specificity - 1) = \max ($TPR - FPR$)



Imbalanced Data



Imbalanced data

Data is imbalanced if there are more records of one class compared to other classes

Imbalanced data may lead to the accuracy paradox

In reality datasets always have some degree of imbalance



Example of imbalanced data

For example: Consider we have information about 500 patients

	Diabetic- Yes	Diabetic-No
Number of records	79	421
% of records	15.8	84.2



Handling imbalanced data

- Up sample minority class
- Down sample majority class
- Change the performance metric
- Try synthetic sampling approach
- Use different algorithm





SMOTE

- SMOTE (Synthetic Minority Oversampling Technique) is one of the most used techniques to deal with an imbalanced data
- It generates the synthetic samples for the minority class (a class with fewer observations)
- Python provides a library that provides different techniques to deal with an imbalanced dataset

```
# install imbalanced-learn using jupyter notebook
# !pip install imbalanced-learn

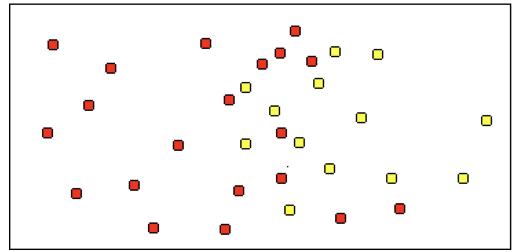
# import SMOTE from imblearn.over_sampling
from imblearn.over sampling import SMOTE
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```



Improper fit

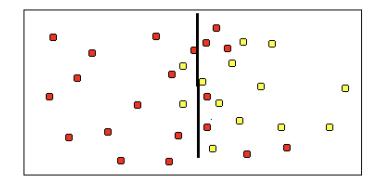
Classify the following data by drawing a line or curve



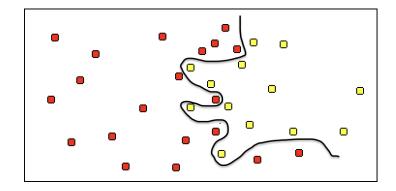


Improper fit

Underfitting

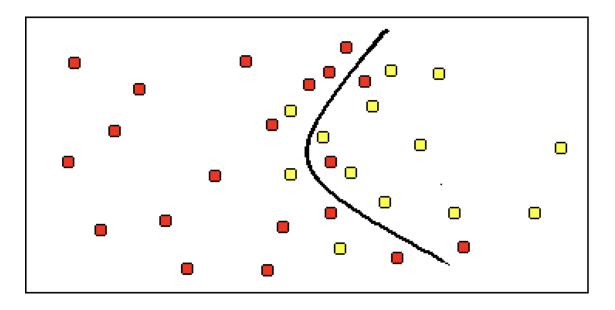


Overfitting





Good fit





Thank You