



Midterm Exam Review

Sravani Vadlamani

Midterm Exam

- In Class on Thursday, March 3, 2022
- Please bring your computer as the exam is on canvas
- Do not forget your calculator

Supervised vs Unsupervised Learning

- Supervised Learning
 - Predict or estimate an output based on one or more inputs
 - Linear regression, logistic regression, boosting, support vector machines
- Unsupervised Learning
 - Learn relationships and structure from data with inputs but no outputs
 - Cluster Analysis

Input vs Output Variables

- Input Variables
 - Independent variables, predictors features
- Output Variables
 - Response or dependent variable
- We believe there is a relationship between Y and at least one of the X's. We model the relation as

$$Y_i = f(X_i) + \varepsilon$$

- Where f is an unknown function and ε is a random error with mean zero
- Parametric vs Non-Parametric Approaches

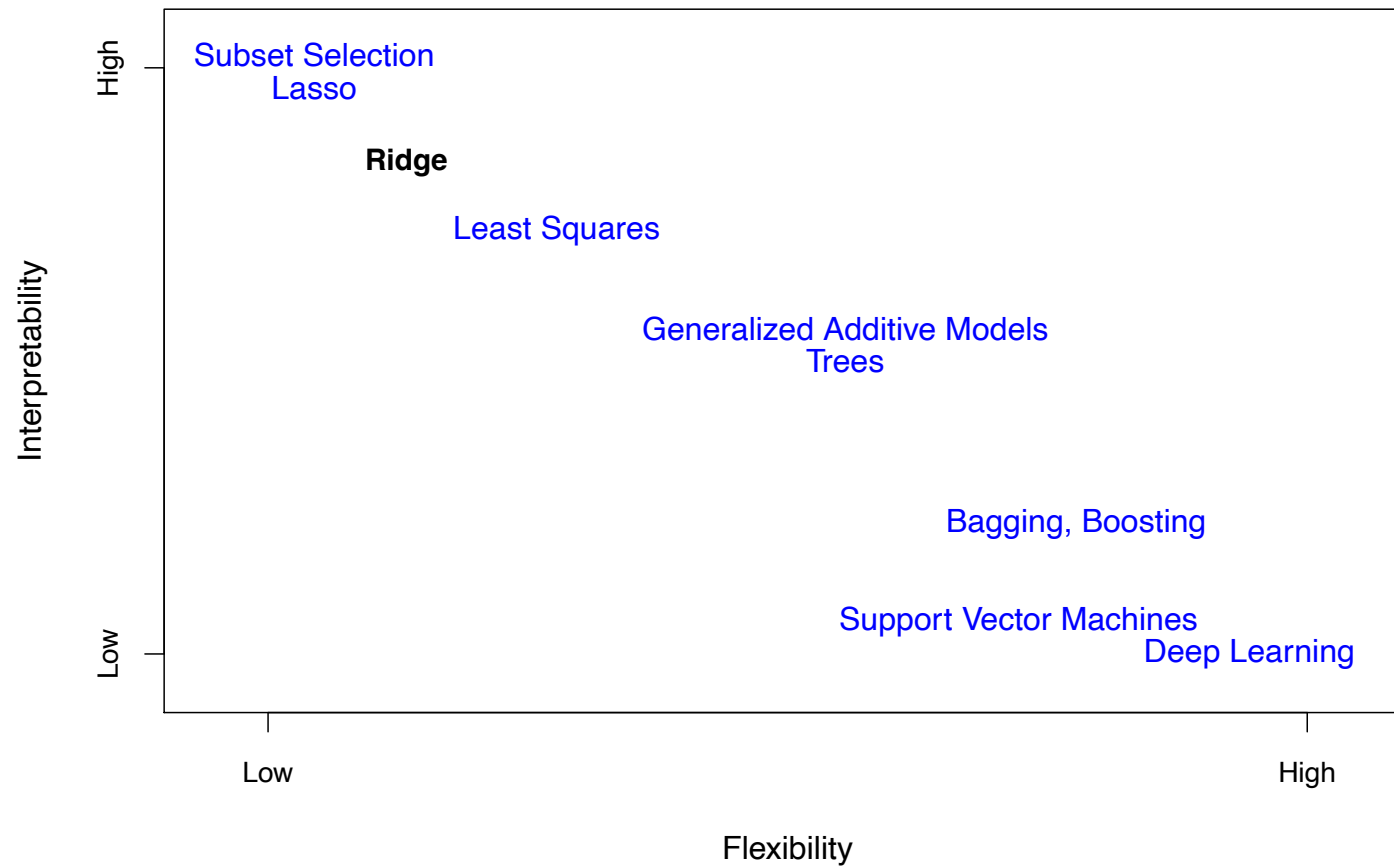
Flexible Models Vs Overfitting

- The more flexible the model, the more realistic it is. However, more flexible models have the disadvantage of requiring a greater number of parameters to be estimated and they are also more susceptible to overfitting.
- Overfitting is a phenomenon where a model closely matches the training data such that it captures too much of the noise or error in the data. This results in a model that fits the training data very well but does not make good predictions under test or in general.

Prediction Accuracy vs Model Interpretability

- Non-linear regression methods are more flexible and can potentially provide more accurate estimates.
- Why not just use a more flexible method if it is more realistic?
- A simple method such as linear regression produces a model which is much easier to interpret (the Inference part is better). For example, in a linear model, β_j is the average increase in Y for a one unit increase in X_j holding all other variables constant.
- Even if you are only interested in prediction, so the first reason is not relevant, it is often possible to get more accurate predictions with a simple, instead of a complicated, model. This seems counter intuitive but has to do with the fact that it is harder to fit a more flexible model.

Prediction Accuracy vs Model Interpretability



Regression vs Classification

- When are both methods used?

Quality of Fit for Regression

- To evaluate the performance of a model, it is necessary to quantify how close the predicted responses are to the observed/actual data
- One common measure of accuracy in regression method is the mean squared error

$$MSE = \frac{1}{n} \sum_{i=1}^n (y_i - \hat{y}_i)^2$$

- Where, \hat{y}_i is the predicted responses

Bias Variance Tradeoff

- Choice of learning method is governed by two competing factors – bias and variance
- Bias refers to the error introduced by modeling a usually extremely complicated problem using a simple problem
- For example – a linear regression model assumes a linear relation between Y and X which may be unlikely in real life thus introducing some bias
- More flexible (or complex) models have less bias

Bias is the error introduced to a complicated problem when modeled with a simple model

Bias Variance Tradeoff

- Variance refers to the amount by which f would change if it were estimated with a different training set
- The more flexible a method is, greater is its variance
- In general, as the flexibility of the statistical method increases, its variance increases and bias decreases
- The relationship between bias, variance, and test set mean squared error is referred to as the bias-variance trade-off. It is called a trade-off because it is a challenge to find a model that has both a low variance and a low squared bias.

Bias Variance Tradeoff

- For any given $x = x_0$, the expected test mean squared error can be decomposed into the sum of the following three quantities:
 - Variance of $f(x_0)$
 - Squared bias of $f(x_0)$
 - Variance of the error term (ϵ)

$$\text{Expected Test MSE} = E(Y - f(x_0))^2 = \text{Bias}^2 + \text{Var} + \underbrace{\sigma^2}_{\text{Irreducible Error}}$$

- To minimize the expected test error, it's necessary to choose a method that achieves both low variance and low bias. It can be seen that the expected test mean squared error can never be less than , the irreducible error.

Linear Algebra Basics

- Matrices
 - Row and Column Vectors
 - Transpose
 - Symmetric Matrix
 - Diagonal Matrix
 - Matrix Addition/Subtraction
 - Matrix Multiplication
 - Hilbert Matrix
 - Trace of a square matrix **Sum of diagonals**

Linear Regression

- Simple and Multiple Linear Regression
- Parameter estimation using least squares
- Forward vs backward variable selection
- Quantitative and Qualitative predictors
- Model Fit
 - Residual Standard Error
 - R-square – proportion of variance explained by the model

Linear Regression

- Issues
 - Non-linearity (if present transform predictor variables)
 - Correlation of error terms
 - Homoscedasticity violation (transform the response variable)
 - Outliers
 - Multicollinearity (VIF) **Variance Inflation Factor**

Classification Techniques

- Logistic regression
- Linear Discriminant Analysis
- Quadratic Discriminant Analysis
- Naïve Bayes
- K-nearest neighbors
- Generalized additive models
- Decision Trees, Random forests, Boosting
- Support Vector Machines

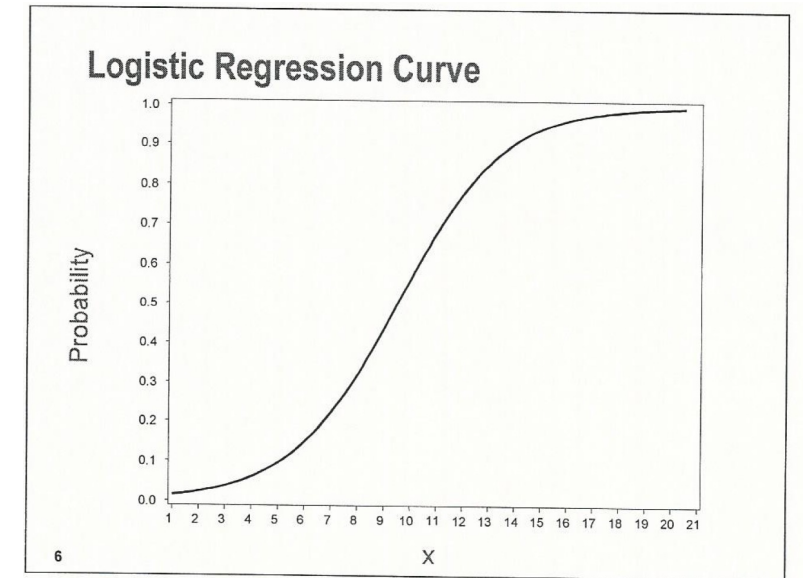
Logistic Regression

- We use a logistic function to model $p(x)$ such that the output is always between 0 and 1 for all values of X

$$p(x) = \frac{e^{\beta_0 + \beta_1 x}}{1 + e^{\beta_0 + \beta_1 x}}$$

- The above equation can be rewritten as

$$\frac{p(x)}{1 - p(x)} = e^{\beta_0 + \beta_1 x}$$



Logistic Regression Parameter Interpretation

- In linear regression, β_1 gives the average change in Y associated with a one-unit increase in X
- In logistic regression, a one-unit change in X yields a β_1 change in the log-odds
 - This is equivalent to multiplying the odds by e^{β_1}
- If β_1 is positive, increasing X will increase $p(x)$
- If β_1 is negative, increasing X will decrease $p(x)$
- The rate of change in $p(x)$ per unit change in X depends on the value of X

Estimating Parameters

- The coefficients β_0 and β_1 must be estimated based on the training
- Logistic regression uses maximum likelihood to estimate β_0 and β_1 such that the predicted probability is as close to the observed classes using the following likelihood function

$$l(\beta_0, \beta_1) = \prod_{i:y_i=1} p(x_i) \prod_{i':y_{i'}=0} (1 - p(x_{i'}))$$

- The estimates β_0 and β_1 maximize the above function

Logistic Regression vs LDA

- In logistic regression we model the conditional distribution of the response variable, given the predictor(s) X i.e., $\Pr(Y=k \mid X = x)$
- In LDA we model the distribution of the predictors separately in each of the response classes Y ($\Pr(X \mid Y=k)$) and then use Bayes theorem to invert the probabilities to estimate the conditional distribution.
- Reasons to prefer LDA over Logistic regression
 - Parameter estimates from logistic regression are surprisingly unstable when there is a substantial separation between classes. Discriminant analysis approaches do not have this issue
 - Logistic regression estimates are less accurate when the predictors are approximately normally distributed, and the sample size (n) is small

LDA with One Predictor - Summary

- In summary, LDA assumes that the observations from each class follow a normal distribution with a class specific mean vector and constant variance across the classes to build a Bayes' theorem-based classifier.

Multiple Predictor LDA

- Assumes $X = (X_1, X_2, \dots, X_p)$ follows a multivariate normal or multivariate Gaussian distribution with a class-specific mean vector and a common covariance matrix.
- Multivariate normal distribution implies that each predictor follows a one-dimensional normal distribution with some correlation between the predictors. The bell shape of the normal distribution will be distorted if the predictors are highly correlated.

Quadratic Discriminant Analysis

- An alternative approach to LDA
- Same assumptions as LDA regarding the observations from each class following a Gaussian/Normal distribution.
- Difference is in the covariance estimation
 - QDA assumes each class has its own covariance matrix.
- This results in assuming that an observation from the k th class follows a distribution of the form $X \sim N(\mu_k, \Sigma_k)$ where Σ_k is the covariance matrix for k th class

Comparison of Methods

- Generally, LDA is better than QDA if there are relatively few training observations and so reducing variance is relevant.
- If the training set is very large that the variance of classifier is not an issue or if the assumption of common covariance matrix is unrealistic, QDA can be a better choice.
- LDA and logistic regression work well when the decision boundary is linear
- QDA gives better results when the decision boundary is moderately non-linear
- K-nearest neighbors (KNN) is a non-parametric approach and outperforms LDA and logistic regression when the decision-boundary is highly non-linear

Metrics

	Condition Positive	Condition Negative
Test Positive	True Positive	False Positive (Type I Error)
Test Negative	False Negative (Type II Error)	True Negative

- Sensitivity = Recall = $P(\text{Test} + \mid \text{Condition} +) = \frac{TP}{TP+FN}$
- Specificity = $P(\text{Test} - \mid \text{Condition} -) = \frac{TN}{FP+TN}$
- False Negative Rate (β) = $P(\text{Test} - \mid \text{Condition} +) = \frac{FN}{TP+FN}$
- False Positive Rate (α) = $P(\text{Test} + \mid \text{Condition} -) = \frac{FP}{FP+TN}$
- Precision = $\frac{TP}{TP+FP}$
- Sensitivity = 1 – False Negative Rate = Power
- Specificity = 1 – False Positive Rate
- F1 Score = $\frac{2 * \text{Precision} * \text{Recall}}{\text{Precision} + \text{Recall}}$
- Accuracy = $\frac{\text{Correct Predictions}}{\text{Total Predictions}}$

Metrics

- Accuracy
 - % of correct predictions
 - One value for the entire model
- Prediction
 - Exactness of the model
 - Each class/label has a value
- Recall
 - Completeness of model
 - Correctly detected over total observations
 - Each class/label has a value
- F1 Score
 - Combines precision and recall (Harmonic mean of precision & recall)
 - Each class/label has a value

ROC Curves

- A curve for simultaneously displaying the positive and negative error types for various thresholds
- ROC stands for “Receiver Operating Characteristics” and derives its name from communications theory (historical)
- The x-axis is False Positive Rate ($=1-\text{sensitivity}$)
- The y-axis is True Positive Rate ($= \text{specificity}$)
- Overall performance is given by the Area Under the Curve, denoted as AUC.
- Larger the AUC, better the classifier i.e., the curve is closer to the top left corner

Linear Model Selection

- Best subset selection
- Stepwise selection
- Model selection criteria
 - C_p , AIC, BIC (lower the better)
 - Adjusted R-square (higher the better)
- Shrinkage methods
 - Ridge Regression (l_2 norm)
 - Lasso Regression (l_1 norm)

Resampling Techniques

- Validation Set Approach
- LOOCV
- K-fold
- Bootstrapping