

unsupervised learning: learning model from unlabeled data
 clustering / dimensionality reduction
 Reinforced learning: improve the performance based on interaction with environment. The feedback is a measure of how well the action was measured by a reward function.

Support Vector Machines
 kernel: mapping to higher-dimensional space
 complexity depends on the number of training samples, not dimensionality
 larger margin - lower error

Performance Matrix
 Actual value

Predicted value	TP	FP
	FN	TN

Decision Tree non
 Linear SVM linear
 SVM with kernel non
 Linear Regression a classifier
 Logistic Regression linear
 Naive Bayes linear

$\phi(w) = w^T w$ is minimized
 Safe margin is allowed errors.

$\phi(w) = w^T w + C \sum \xi_i$
 penalty $C \rightarrow 0$ underfitting
 $C \rightarrow \infty$ overfitting

SVM Weakness:

- ① sensitive to noise
- ② Standard SVM only consider 2 classes
 ↓ build multiple SVMs
- ③ select a specific kernel and parameters is usually done by see and try

Linear Regression: Least Square Fitting

minimize: $\sum_{i=1}^n [y_i - (\beta_0 + \beta_1 x_i)]^2$
 $S_{xy} = \sum_{i=1}^n x_i y_i - \frac{1}{n} (\sum_{i=1}^n x_i) (\sum_{i=1}^n y_i)$
 $S_{xx} = \sum_{i=1}^n x_i^2 - \frac{1}{n} (\sum_{i=1}^n x_i)^2$
 $S_{yy} = \sum_{i=1}^n y_i^2 - \frac{1}{n} (\sum_{i=1}^n y_i)^2$
 $\hat{\beta}_0 = \bar{y} - \hat{\beta}_1 \bar{x}$
 $\hat{\beta}_1 = \frac{S_{xy}}{S_{xx}}$
 Multiple: $\hat{\beta} = (X^T X)^{-1} X^T Y$

Naive Bayes \rightarrow MAP

$\arg \max_y P(y|x) = \arg \max_y P(y) \prod_{i=1}^n P(x_i|y)$

Logistic Regression directly compare $P(y|x)$
 Naive Bayes use Bayes Theorem to compare $P(y|x)$
 NB + Gaussian Basis Function
 \propto LR + sigmoid

Logistic Regression model the $P(y|x)$ as a logistic function. The logistic is a weighted linear combination of the features. Linear regression itself is based on linear feature function to regress.

\rightarrow Or using total probability:

$P(S) = P(S|yes)P(yes) + P(S|no)P(no) = (3/4)(1/4) + (3/5)(3/4) = \frac{5}{8}$

NB: $P(y|S, w) = P(S, w|yes)P(yes)/P(S, w) = \frac{10}{31}$

NB: $P(S, w|yes) = P(S|yes)P(w|yes) = \frac{6}{31}$

$\rightarrow P(S, w) = P(S, w|yes) + P(S, w|no)P(no)$
 $= P(S|yes)P(w|yes)P(yes) + P(S|no)P(w|no)P(no)$
 $= \frac{31}{210}$

NB: input features x_i is independent given label Y

LR: Least square is not suitable, Maximum likelihood estimation instead

(TPR) Recall: $\frac{TP}{FN+TP}$

when false negatives is catastrophic, e.g. disease detection

(PPV) Precision: $\frac{TP}{TP+FP}$
 NPV = $\frac{TN}{TN+FN}$

when being right (positive prediction) is correct/outweighs detecting all positives, e.g. recommendation system

① Training set:

↓

② Validation set: hyperparameter tuning and model selection

↓

③ Test data

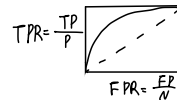
K-fold cross validation is used when we have little data



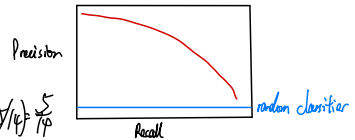
F1-score: $2 * \frac{\text{Precision} * \text{Recall}}{\text{Precision} + \text{Recall}}$

(TNR) Specificity = $\frac{TN}{TN+FP}$

ROC Curves: the trade-off between TP rate and FP rate



PR Curve: between TP rate and positive predictive value



Oscar's Razor

Given 2 models with similar generalization errors, the simple model is preferred.

Probabilistic classifiers: yields a probability distribution. includes: Random forest, NB, LR

Deterministic classifiers: no model \rightarrow specific feature space includes: Decision Tree, SVM

Generative classifiers: learn $P(x, y)$, calculate $P(y|x)$ to find max $P(y)$ NB, Bayes Naive, Hidden Markov

Discriminative classifiers: learn $P(y|x)$ directly LR, SVM

Decision Trees:

when have N attributes, $2^{(2^N)}$ trees

Choose feature $X_i = H(Y|X_i)$ ↓

$I(X; Y) = H(Y) - H(Y|X)$ ↑

$H(X) = -\sum p(x) \log_2 p(x)$

$H(Y|X) = \sum p(x) H(Y|X=x)$

$I(X; Y) = \sum \sum P_{x,y}(x,y) \log \left(\frac{P_{x,y}(x,y)}{P_x(x)P_y(y)} \right)$

Gini = $1 - \sum P_i^2$

classification Error = $1 - \max P_i$

Decision Tree Depth ↑ \rightarrow overfitting

Ensemble learning - Random Forest

reduce overfitting and variance without decreasing

performance

Bagging - Bootstrap Aggregating

bootstrapping: Random sampling with replacement

train multiple decision trees & search all features

to split on for each tree

Aggregating: Combine multiple predictions via averaging or majority vote

NN: supervised learning

perceptron: $y = \sigma(b + \sum_{i=1}^m w_i x_i) = \{0, 1\}$

Multi-layer perceptron: $y_j = \sigma(\sum_{i=1}^m w_{ij} x_i + b_j)$

↳ feed forward network

Common Activation Functions:

Sigmoid/tanh/ReLU/Leaky ReLU

Maxout/ELU



NN = Function Approximation

Feedforward NN: no loops input → hidden layers → output

Recurrent: use feedback

RNN: ① employ feedback
not necessarily stable

② forecasting time series data
language translation

Hopfield Networks: fully connected

CNN → Image

Transformer Networks: Natural language Processing

Generative Adversarial Network:

Single perceptron can find linear max margin ✓
LR solution ✓

NB → inference in $P(\text{labels} | \text{features})$

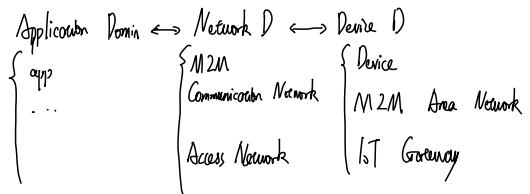
$$B: P(S | (F_1, F_2)) = P(F_1, F_2 | S) P(S) / P(F_1, F_2)$$

$$\text{NB: } P(S | (F_1, F_2)) = \frac{P(F_1, F_2 | S) \times P(S)}{P(F_1 | S) \times P(F_2 | S) \times P(S)}$$

NN optimization: ① Gradient Descent
② Backpropagation

Computing & Intelligence drive IoT

IoT Architecture:



↑ Bottom End: Basic, Resource constrained, *environment monitor*

In the Middle: support localization, full protocol, *in home, industrial*

↓ Top End: *Utility / Medical use*

Technical Issue: ① Naming, Addressing, Routing

② Scalability for network architecture

③ Power saving & management

④ Security

Challenges: ① Sensitive Latency Requirements

② Network Bandwidth Constraint

③ Resource Constrained Device