

Applied AI & ML Industry Projects Lab (AAM-IPL)

Transforming Learning into Industry Solutions



Week-5

SVM Overview and Project Announcement

V Semester - ML and AIUP, Aug-Oct 2024

Session Date and Time: 19th Oct 2024 10:30 AM IST - 12:00 Noon IST

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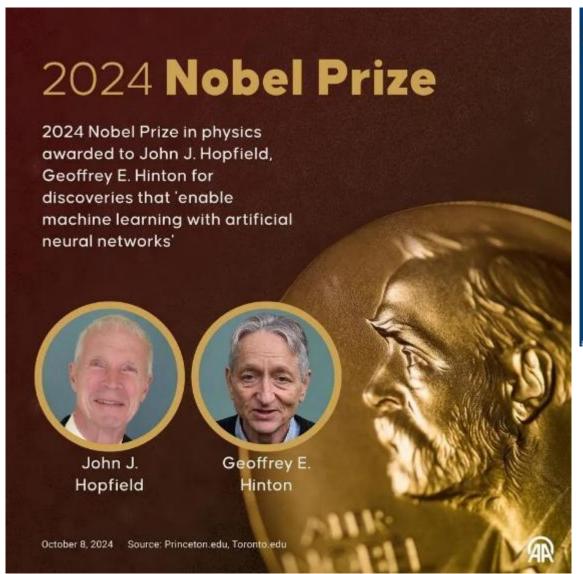
Thank you for your understanding and cooperation.

Lecture Outline



- Classification
- Real-world Classification Use Cases
- Classification Algorithms and Comparison
- Linear Classifiers
- Maximum Margin Classifier
- Constrained Optimisation Lagrangian Duality
- Primal and Dual SVM
- Non-linear Data and Higher Dimensional Classification
- Kernel Tricks and Functions
- Common Kernel Functions
- Kernel SVM and Decision Boundaries
- Kernel Tricks Vs PCA
- SVM with Slack Variables
- Project Demo
- Q&A

Al Steals The Spotlight at 2024 Nobel Prizes





David Baker

University of Washington, Seattle, WA, USA Howard Hughes Medical Institute, USA "for computational protein design" and the other half jointly to

Demis Hassabis
Google DeepMind, London, UK
John M. Jumper
Google DeepMind, London, UK
"for protein structure prediction"

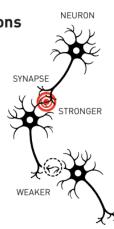
Source: Press release: 8th Oct 2024 - The Nobel Prize in Physics 2024 - Nobel Prize.org



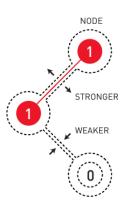
Al Steals The Spotlight at 2024 Nobel Prizes

Natural and artificial neurons

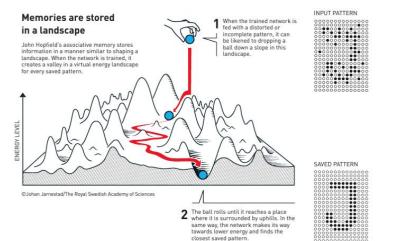
The brain's neural network is built from living cells, neurons, with advanced internal machinery. They can send signals to each other through the synapses. When we learn things, the connections between some neurons gets stronger, while others get weaker.

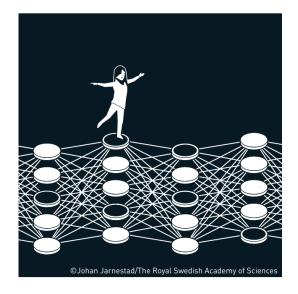


Artificial neural networks are built from nodes that are coded with a value. The nodes are connected to each other and, when the network is trained, the connections between nodes that are active at the same time get stronger, otherwise they get weaker.

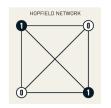


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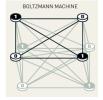




Different types of network

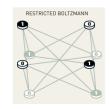


John Hopfield's associative memory is built so that all the nodes are connected to each other. Information is fed in and read out from all the nodes.



Visible Hidde nodes

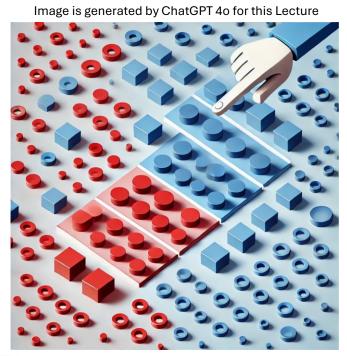
Geoffrey Hinton's Boltzmann machine is often constructed in two layers, where information is fed in and read out using a layer of visible nodes. They are connected to hidden nodes, which affect how the network functions in its entirety.



In a restricted Boltzmann machine, there are no connections between nodes in the same layer. The machines are frequently used in a chain, one after the other. After training the first restricted Boltzmann machine, the content of the hidden nodes is used to train the next machine, and so on.

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Classification



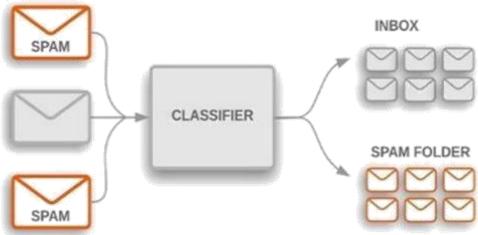


Image Courtesy: (1) Classification Algorithms in Machine Learning. | LinkedIn

- Classification is the process of assigning data into classes or categories - predefined or otherwise, based on its features
- Determines to which class an observed data belongs to
- The decision boundary separates two distinct groups, each representing a different class
- In the image on to the left, a classifier distinguished red circles and squares belonging to one class from blue circles and squares belonging to another
- Support Vector Machines SVM and other classification algorithms learn to create such boundaries for accurate data class predictions
- The goal of classification is to make decisions based on these separated groups
 - Identifying a received email as spam
 - Marking a credit card transaction as fraud
 - Defining customer categories

Real-World Use Cases of Classification

Decision Making and Prediction

- Allows us to predict categorical outcomes and make informed decisions based on historical data
- Example: Healthcare Predict whether a patient has a disease or not based on medical data

Tasks Automation

- Automation of sorting and categorizing data which otherwise would be laborious and error-prone when done by humans
- Example: Spam classifier/detector in email server to automatically sort and filter each received emails into span and genuine emails

Understanding Patterns and Trends

- Helps us uncover hidden patterns in data that would otherwise not be immediately obvious through manual inspection
- Example: Online retail companies use classification to understand customer behaviour to identify whether a customer is likely to purchase a product based on their browsing history

Real-World Use Cases of Classification

Improved Decision Making in Business

- Companies use classification to segment customers, predict customer churn, and to target marketing efforts more effectively for growing business
- Example: A bank can use classification to determine which customers are likely to default on a loan, allowing them to mitigate risks in loan repayment

Handling Complex High-Dimensional Data

- Advanced classification algorithms like SVM or deep learning, can find patterns in complex, highdimensional datasets that humans, otherwise, would struggle to analyze
- Example: In facial recognition or autonomous vehicle driving, classification algorithms help identify objects or people in images or in live video

NLP

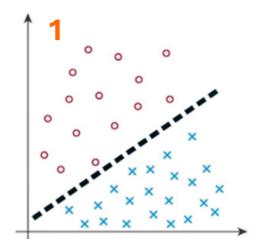
Sentiment analysis and text classification

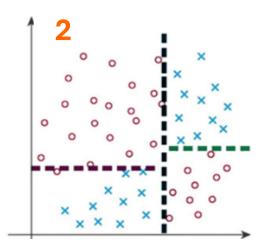
Classification Algorithms

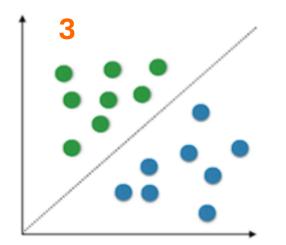
- Based on Output
 - Binary Classifiers 2 output classes
 - Logistic Regression, Simple Naïve Bayes, etc.
 - Multinomial Classifiers More than 2 output classes
 - Multi-monomial Logistic Regression/ Softmax, Non-linear Naïve Bayes, KNN, Decision Tree Classifier, etc.
- Based on Data Characteristics
 - Linear Classifier Classify data using a linear decision boundary a straight line or hyperplane
 - Non-Linear Classifier Classifying using complex, non-linear decision boundaries

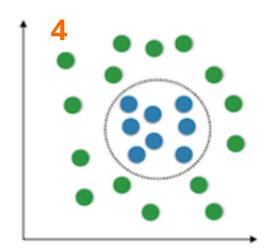


Classification Algorithms - Linear and Non-linear Classifiers









Linear Classifiers

- Examples: Logistic Regression, SVM, Linear Discriminant Analysis, etc.
- Best for
- Data that can be separated by a straight line (linearly separable data).
- Advantages
- Simpler and faster to compute
- Easier to interpret
- Work well for high-dimensional, sparse data (e.g., text classification)
- Limitations
- Cannot handle complex, non-linear patterns in data

Non-linear Classifiers

- Example : Kernel SVM, DTC, Random Forest, KNN, Neural Networks, etc.
- Best for
 - Data with complex, non-linear relationships
- Advantages
 - Can model complex and intricate patterns
 - More flexible, able to handle diverse types of data
- Limitations
 - Often computationally more expensive
 - May require more data to train
- More challenging to interpret

Key Comparison

Decision Boundary

• Linear classifiers use straight lines or hyperplanes, while non-linear classifiers can create curves or complex boundaries

Performance

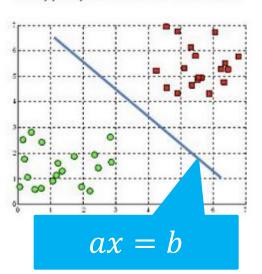
• Non-linear classifiers generally perform better when data isn't linearly separable

Computational Cost

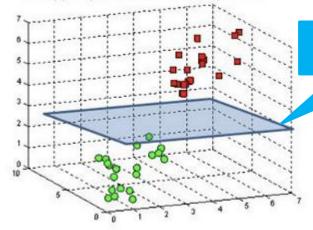
- Linear classifiers are faster and simpler
- Non-linear ones are more computationally expensive but powerful for complex tasks

Classifier Decision Boundaries

A hyperplane in \mathbb{R}^2 is a line



A hyperplane in \mathbb{R}^3 is a plane

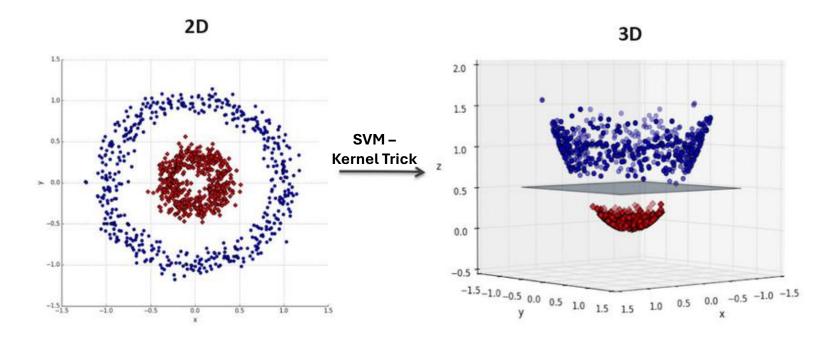


 $\bar{a}^T\bar{x}=b$

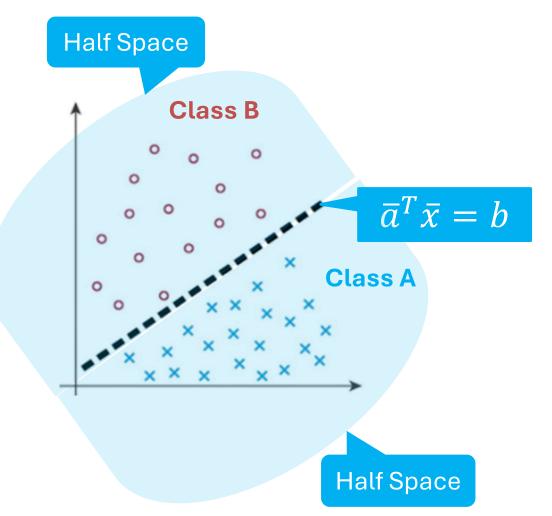
2-D: $a_1x_1 + a_2x_2 = b$

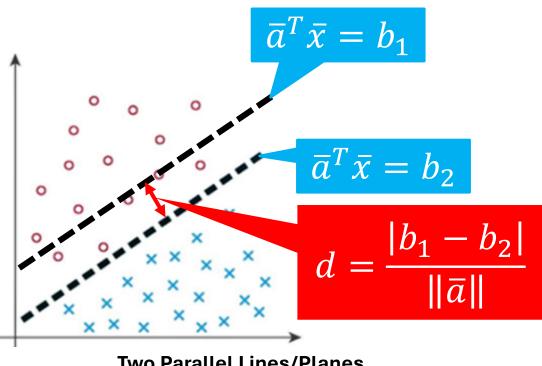
3-D: $a_1x_1 + a_2x_2 + a_3x_3 = b$

 $n-D: a_1x_1 + a_2x_2 + \cdots + a_nx_n = b$



Linear Classifier





Two Parallel Lines/Planes

General Structure of a Linear Classifier:

Class A: \mathcal{C}_A : $\overline{a}^T \overline{x} \leq b$

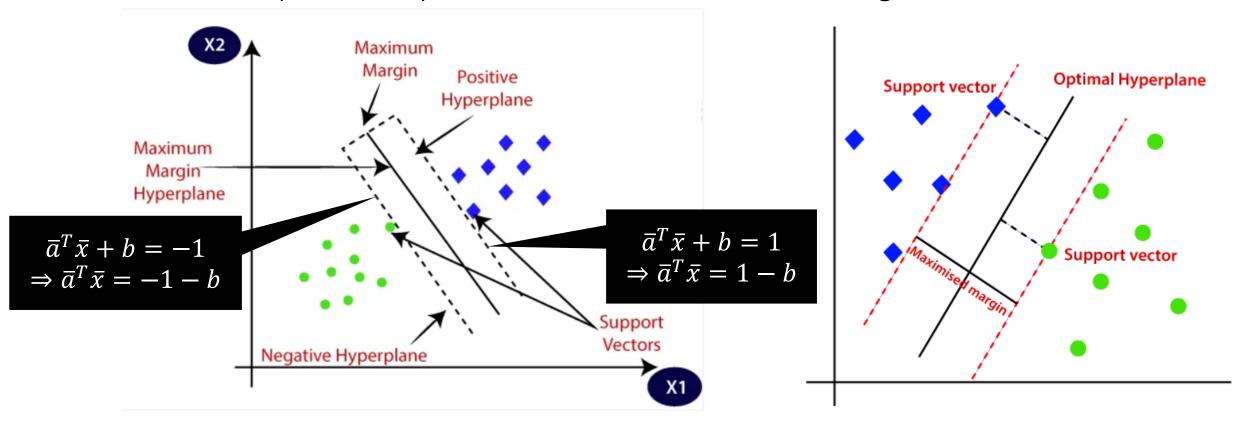
Class B: C_R : $\overline{a}^T \overline{x} > b$

Distance Between Two Parallel Lines:

$$d = \frac{|b_1 - b_2|}{\sqrt{a_1^2 + a_2^2 + \cdots + a_n^2}}$$

Support Vector Machines – Maximum Margin Classifier

• Objective of SVM algorithm is to find **an optimal hyperplane/decision boundary** in an n-dimensional data space that separates classes with the **maximum margin**

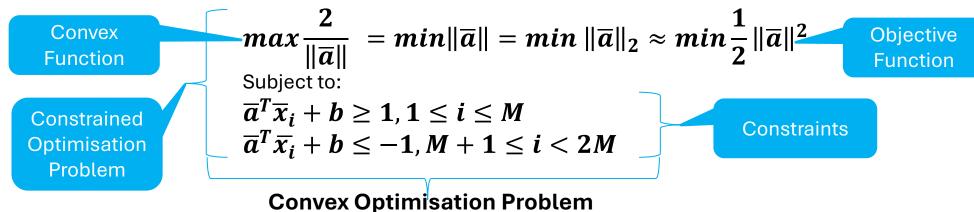


Margin – Distance between the positive and negative hyperplanes:

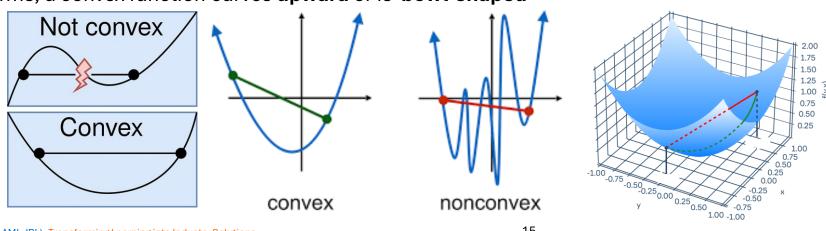
$$d = \frac{|1 - b - (-1 - b)|}{\|\overline{a}\|} = \frac{|2|}{\|\overline{a}\|}$$

Linear Maximum Margin Classifier - SVM Classifier

• Maximize the margin/distance (d) between the positive and negative hyperplanes for all data points in training set 1 ... 2*M*



- Convex Function
 - The line segment connecting any two points on the function graph, lies above or on the graph of the function
 - f(x) is convex, if for any two points \bar{x}_1 and \bar{x}_2 in its domain and for any $\lambda \in [0,1]$, the inequality $f(\lambda \bar{x}_1 + (1-\lambda)\bar{x}_2) \le 1$ $\lambda f(\bar{x}_1) + (1-\lambda)f(\bar{x}_2)$
 - $\bullet \ \forall \ \bar{x}_1, \bar{x}_2 \in \mathcal{D} \ and \ \lambda \in [0,1] \Rightarrow f(\lambda \bar{x}_1 + (1-\lambda)\bar{x}_2) \leq \lambda f(\bar{x}_1) + (1-\lambda)f(\bar{x}_2)$
 - In simpler terms, a convex function curves upward or is bowl-shaped



Constrained Optimisation – Lagrangian Duality

Standard optimisation problem form - aka Primal Problem

```
p^* = min f(x)
Subject to (Constraints):
g_i(x) \le 0, i = 1,2, \dots m – Inequality constraints
h_i(x) \ge 0, j = 1,2, \dots p – Inequality constraints
c_k(x) = 0, k = 1,2, \dots t – Equality constraints
x \in \mathbb{R}^n, domain \mathcal{D}, optimal value p^*
```

- Lagrangian Duality aka Dual Optimisation Problem
 - A simplified Lagrange Function aka Lagrangian is formulated:
 - By combining objective function and constraints using Lagrange multipliers (λ, μ, ν)
 - This mathematical formulation reflects the **impact of constraints** in the solution
 - $\mathcal{L}(x,\lambda,\mu,\nu)=f(x)+\sum_{i=1}^m\lambda_ig_i(x)-\sum_{j=1}^p\mu_jh_j(x)+\sum_{k=1}^t\nu_kc_k(x)$ where $\lambda_i\geq 0$, $\mu_j\leq 0$, ν_k no constraints
 - Minimize the **Lagrangian** with respect to the original/primal variable x while treating the Lagrange multipliers as fixed
 - $\Phi(x,\lambda,\mu,\nu) = \inf / \min \mathcal{L}(x,\lambda,\mu,\nu)$
 - Result is **dual function** $\phi(\lambda, \mu, \nu)$ and depends only on Lagrange multipliers λ, μ, ν
 - A dual problem is constructed to maximise the dual function w.r.t to Lagrange multipliers subject to the feasibility conditions on Lagrange multipliers
 - $\blacksquare \max_{\lambda \ge 0, \mu \le 0, \nu} \phi(\lambda, \mu, \nu)$
- When the maximum margin classifier is solved using Lagrangian Duality ⇒ Dual SVM

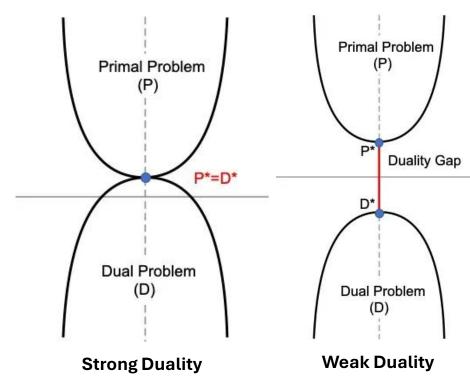
Primal SVM and Dual SVM

Primal SVM

- Directly solves for the hyperplane/decision boundary in feature space
- Suitable for low-dimensional problems
- The maximum margin classifier is solved using primal optimisation problem

Dual SVM

- Preferred when the number of features is larger than the number of samples
- Instead of solving for the weights \overline{a} and b of the hyperplane directly from **primal optimisation problem**, dual SVM solves for the **Lagrange multipliers** making it more flexible for non-linear decision boundaries
- Reduced size of the problem from a high-dimensional feature space to a lower-dimensional dual space
- The maximum margin classifier is solved using dual optimisation problem



Dual SVM Solution

• Solve $minrac{1}{2}\|\overline{a}\|^2$

Subject to:

$$\overline{a}^T \overline{x}_i + b \ge 1, 1 \le i \le M$$

 $\overline{a}^T \overline{x}_i + b \le -1, M + 1 \le i < 2M$

Two constraints can be combined into one

$$y_i(\bar{a}^T\bar{x}_i+b)\geq 1, \forall i$$

Where:

$$y_i = \begin{cases} +1, & i = 1,2, \dots M \\ -1, & i = M+1, \dots 2M \end{cases}$$

Lagrangian

$$\mathcal{L}(\bar{a}, b, \lambda) = \frac{1}{2} \|\bar{a}\|^2 - \sum_{i=1}^{2M} \lambda_i [y_i(\bar{a}^T \bar{x}_i + b) - 1] = \frac{1}{2} \bar{a}^T \bar{a} - \sum_{i=1}^{2M} \lambda_i [y_i(\bar{a}^T \bar{x}_i + b) - 1]$$

- Solving for \bar{a} , b minimising $\mathcal{L}(\bar{a},b,\lambda)$
 - $\phi(\lambda, \bar{x}) = \inf / \min_{x} \mathcal{L}(\lambda)$
 - With respect to a
 - $\bullet \frac{\partial \mathcal{L}}{\partial a} = \nabla_{\bar{a}}(\mathcal{L}(\bar{a}, b, \lambda)) = \bar{a} \sum_{i=1}^{2M} \lambda_i y_i x_i = 0 \Rightarrow \bar{a} = \sum_{i=1}^{2M} \lambda_i y_i x_i$
 - The points for which $\lambda_i \neq 0$ are called **Support Vectors**
 - With respect to b

• Substituting \bar{a} and b in Lagrangian and simplifying for **dual function** $\phi(\bar{x}, \lambda) = \sum_{i=1}^{2M} \lambda_i - \frac{1}{2} \sum_{i,j=1}^{2M} \lambda_i \lambda_j y_i y_j \bar{x}_i^T \bar{x}_j$

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Dual SVM Solution contd...

- Therefore dual problem $\max_{\lambda \geq 0} \phi(\lambda, \overline{x})$ can be formulated as:
 - $max \left[\sum_{i=1}^{2M} \lambda_i \frac{1}{2} \sum_{i,j=1}^{2M} \lambda_i \lambda_j y_i y_j \overline{x}_i^T \overline{x}_j \right]$

Subject to:

$$\lambda_i \geq 0, \forall i$$

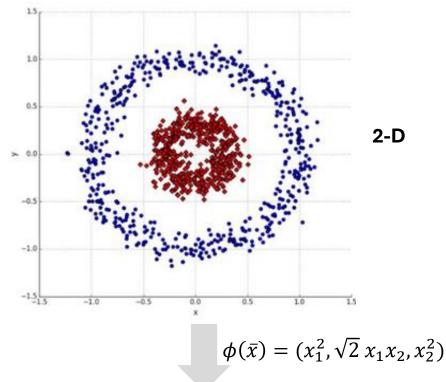
$$\sum_{i=1}^{2M} \lambda_i = 0$$

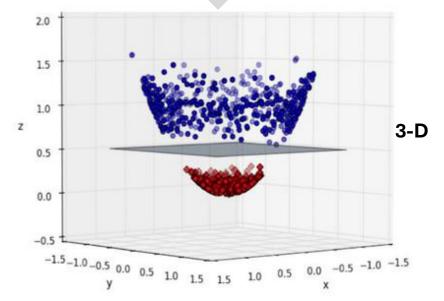
- Dual problem is a quadratic programming problem objective is quadratic and constraints are linear
- Solved for λ_i using standard quadratic solvers
 - Python and SVM solvers cvxopt, libsvm, sklearn.svm etc.
 - \bar{a} and b are computed as below
 - $\overline{a} = \sum_{i=1}^{2M} \lambda_i y_i \overline{x}_i$
 - b is found by using any one support vector $(\lambda_i > 0)$ and solving $y_i(\overline{a}^T \overline{x}_i + b) = 1$
- New data point classification
 - $f(\bar{x},b) = \bar{a}^T \bar{x} + b > 0 \Rightarrow$ data point is classified as y = 1 Class 1
 - $f(\bar{x},b) = \bar{a}^T\bar{x} + b < 0 \Rightarrow$ data point is classified as y = -1 Class 0
- In its simplest form, SVM works well for linearly separable data
- In real-world, data is often **not linearly separable** in the original input space

How can we achieve linear separability for a non-linearly separable real world data?

Non-linear Data and Higher Dimensional Classification

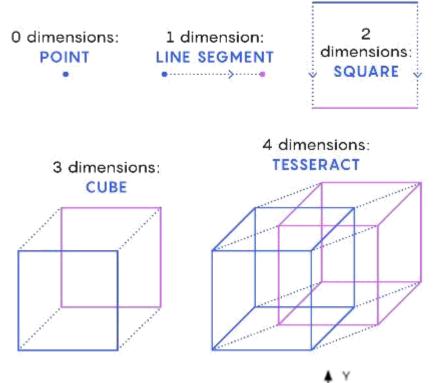
- Dual problem
 - $\max\left[\sum_{i=1}^{2M}\lambda_i-\frac{1}{2}\sum_{i,j=1}^{2M}\lambda_i\lambda_jy_iy_j\overline{x}_i^T\overline{x}_j\right]$ Subject to: $\lambda_i\geq 0$, $\forall i$ and $\sum_{i=1}^{2M}\lambda_i$ =0
- Data points \bar{x}_i , \bar{x}_j appear in the form of inner dot product $\bar{x}_i^T \bar{x}_j$ dot/inner product $\langle \bar{x}_i, \bar{x}_j \rangle$
- A non-linear data set in its current dimensional space χ might become linearly separable if transformed into a higher-dimensional space ν
 - Set of 2-D points forming concentric circles non-linear data set
- The input dataset is transformed into a higher dimensional space using **mapping function** $\phi(\overline{x})$
 - Example : 2-D point (x_1, x_2) is transformed into 3-D space using $\phi(\bar{x}) = (x_1^2, \sqrt{2} x_1 x_2, x_2^2)$
 - Possible to separate classes linearly with a hyperplane in 3-D
- Explicit computation of mapped vectors and their inner product $\phi(\bar{x}_i)$. $\phi(\bar{x}_j)$ in higher dimensional feature space becomes computationally expensive

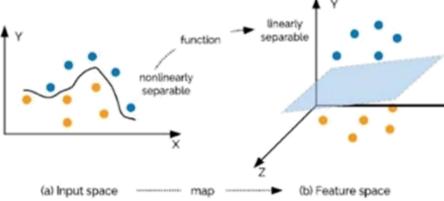




Kernel Tricks/Methods

- Kernel Tricks or Methods owe their name to Kernel Functions
- These kernel functions enable to operate in a high-dimensional, implicit feature space without needing to compute the higher dimensional space coordinates but simply computing the inner products between the images of all pairs of data in the feature space
- For all $\bar{x_i}$, $\bar{x_j}$ in the input space χ , certain functions $K(\bar{x_i}, \bar{x_j})$ can be expressed as an inner product in another/higher dimensional space ν
 - $K: \chi \times \chi \to \mathbb{R}$ is often referred to as a Kernel of Kernel function
- The dot product computation in the higher dimensional space ν is made much simpler using
 - $\phi: \chi \to \nu$ satisfying $K(\bar{x}_i, \bar{x}_j) = \langle \phi(\bar{x}_i), \phi(\bar{x}_j) \rangle_{\nu}$
 - An explicit representation for ϕ is NOT necessary as long as ν is an inner product
- Kernel Function $K(\overline{x}_i, \overline{x}_j)$ that produces equivalent dot product in transformed higher dimensional space $\phi(\overline{x}_i)^T$. $\phi(\overline{x}_j)$
- This trick helps avoid the explicit computation of mapped vectors and their inner product
- Kernel function $K(\bar{x}_i, \bar{x}_j)$ is a generalisation of the dot product in higher feature space $\phi(\bar{x})$





Common Kernel Functions

Linear Kernel

- $K(\overline{x}_i, \overline{x}_j) = \overline{x}_i \cdot \overline{x}_j$
- Equivalent of no transformation works for linearly separable data

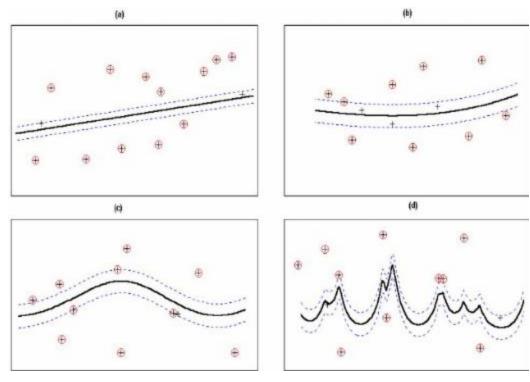
Polynomial Kernel

- $K(\overline{x}_i, \overline{x}_j) = (\overline{x}_i, \overline{x}_j + c)^d$
- Maps data into polynomial feature space to model polynomial decision boundary in original space

Radial Basis Function aka Gaussian Kernel

•
$$K(\overline{x}_i, \overline{x}_j) = e^{\left(-\frac{\left\|\overline{x}_i - \overline{x}_j\right\|^2}{2\sigma^2}\right)}$$

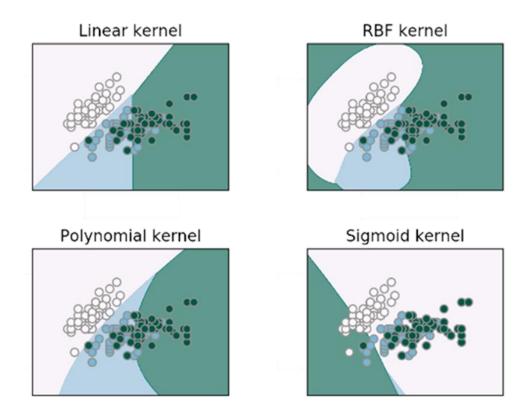
- Maps data into an infinite-dimensional feature space enabling highly flexible decision boundary
- A right kernel function is chosen based on data complexity and distribution



(a) Linear, (b) Polynomial, (c) Gaussian RBF and (d) Exponential RBF

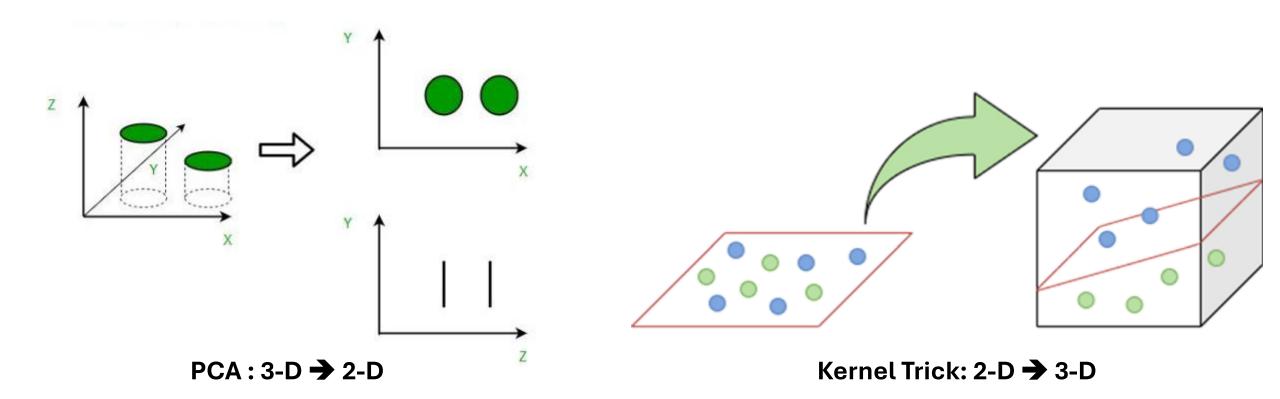
Kernel SVM and Decision Boundaries

- Substituting the Kernel Function into the dual problem
 - $max \left[\sum_{i=1}^{2M} \lambda_i \frac{1}{2} \sum_{i,j=1}^{2M} \lambda_i \lambda_j y_i y_j \overline{x_i^T} \overline{x_j} \right] \Rightarrow max \left[\sum_{i=1}^{2M} \lambda_i \frac{1}{2} \sum_{i,j=1}^{2M} \lambda_i \lambda_j y_i y_j \overline{K}(\overline{x_i}, \overline{x_j}) \right]$
 - Subject to: $\lambda_i \geq 0$, $\forall i$ and $\sum_{i=1}^{2M} \lambda_i = 0$
- The decision boundary after incorporating the Kernel Function
 - $f(\overline{x}, b) = sign(\sum_{i} \lambda_{i} y_{i} K(\overline{x}_{i}, \overline{x}_{i}) + b)$



Kernel Tricks Vs PCA

- PCA reduces dimensionality, Kernel Trick effectively increases dimensionality
- **PCA** performs **linear** operations on the original feature space to find principal/lower dimensional components
- Kernel Trick applies a non-linear transformation via kernel functions to implicitly increase dimensionality
- PCA simplifies, while Kernel Trick enables richer separation

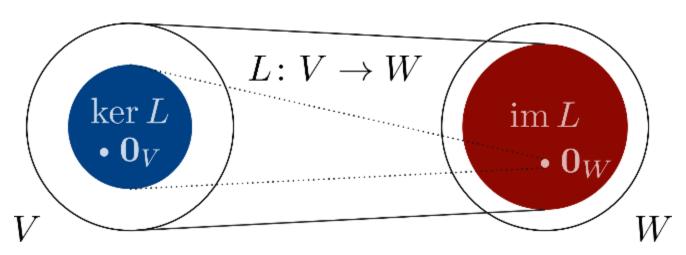


Sidebar Topic - Kernels in Linear Algebra

- Kernel of a linear map/matrix aka null space
- Linear map $L: V \to W V$, W are two vector spaces
- $ker(L) = \{v \in V \mid L(v) = 0\} = L^{-1}(0)$
- L is a linear sub-space of domain V
- Two elements of V have the same image in W iff their difference lies in the kernel of L
 - $v_1, v_2 \in V, L(v_1) = L(v_2) \iff (v_1 v_2) \in V$
 - $\Rightarrow L(v_1) L(v_2) = L(v_1 v_2) = 0$
- Rank-nullity theorem (V is finite dimensional)
 - dim(ker L) + dim(image L) = dim(V)

Linear Map/Transform

$$Lx \iff \begin{bmatrix} l_{11}x_1 + l_{12}x_2 + \cdots l_{1n}x_n \\ l_{21}x_1 + l_{22}x_2 + \cdots l_{2n}x_n \\ \vdots \\ l_{m1}x_1 + l_{m2}x_2 + \cdots l_{mn}x_n \end{bmatrix}_{m \times n}$$

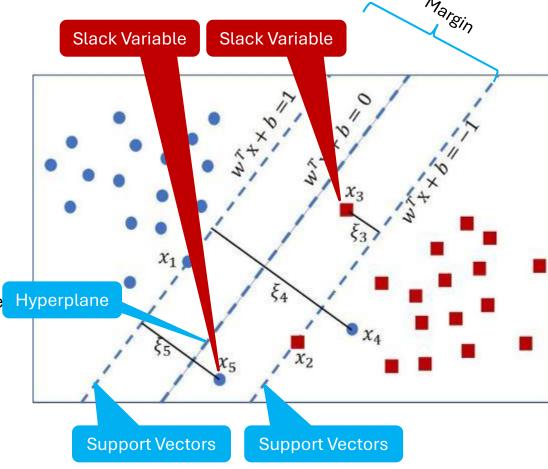


SVM With Slack Variables

- Primal SVM Classifier aka Hard Margin SVM works well when the data is perfectly linearly separable
- Real-world data often contains **Noise** and/or **overlaps** between classes
- One way to handle **noise** and **overlaps** is using **slack variables** (ξ_i) to allow some data points to lie inside the margin or on the wrong side \rightarrow **Soft Margin SVM**
- Optimisation problem with slack variables
 - $\min_{\overline{a},b,\xi} \left[\frac{1}{2} \|\overline{a}\|^2 + C \sum_{i=1}^{2M} \xi_i \right]$

Subject to:

- $y_i(\overline{a}^T\overline{x}_i+b) \geq 1-\xi_i, \ \xi_i \geq 0 \ \forall i$
- $\xi_i \geq 0$ the point is either inside the margin or mis-classified
- C Regularisation parameter to control the trade-off between maximising the margin and minimizing the slack
- Purpose of Slack Variables
 - Classification flexibility works for both linearly separable and nonseparable data
 - Trade-off Control: C allows balancing a larger margin and fewer violations
- Key Points
 - Slack Variables ξ_i allow some points to be inside the margin or misclassified
 - Soft Margin SVM balances maximizing margin and minimizing classification errors using ${\cal C}$
 - SVM handle noisy, overlapping, or non-linearly separable data with slack variables
 - Soft Margin SVM is more flexible, generalizable model



Accuracy and Efficiency

- Accuracy
 - Accuracy represents the proportion of correctly classified samples (both true positives and true negatives) out of the total number of samples.
 - $Accuracy = \frac{TP+TN}{TP+TN+FP+FN}$
- Precision
 - Precision measures how many of the samples predicted as positive are actually positive. It focuses on the quality of the positive predictions.
 - $Precision = \frac{TP}{TP+FP}$
- Recall aka Sensitivity aka TPR
 - Recall measures how many of the actual positive samples were correctly classified. It focuses on the ability to find all positive instances
 - $Recall = \frac{TP}{TP + FN}$
- F1-Score
 - The F1-score is the harmonic mean of precision and recall. It is useful when the dataset has an uneven class distribution.
 - $F1 Score = 2 \times \frac{Precision \times Recall}{Precision + Recall}$
- Confusion Matrix
 - A confusion matrix is a table used to describe the performance of a classification model. It provides insight into the following
 - True Positives (TP): Correctly predicted positive instances
 - True Negatives (TN): Correctly predicted negative instances
 - False Positives (FP) (Type I Error): Incorrectly predicted positive instances
 - False Negatives (FN) (Type II Error): Incorrectly predicted negative instances
- RoC Curve
 - The ROC curve is a graphical representation of a classification model's performance across all classification thresholds. It plots the True Positive Rate (Recall) against the False Positive Rate.
- Feature Importance
 - Relative contribution of each feature to the prediction made by a model. It provides insight into which features have the most influence on the target prediction.

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Key Takeaways

- SVM is powerful for classification
- Works well in high-dimensional spaces and for non-linearly separable data with the kernel trick
- Effective for imbalanced data with hyperparameter tuning

Project Overview

Project Title

Diagnosing breast cancer (malignant or benign tumor) based on the diagnostic data

Project Objective

- This project implements a Support Vector Machine (SVM) classifier to predict the class of breast cancer (breast cancer detected or not detected)) from the provided dataset.
- The dataset is loaded from sklearn
- The breast cancer dataset consists of 569 samples, each representing a patient with a set of features.
- The dataset is used for training a SVM classifier and for evaluating performance metrics such as accuracy, precision, recall, F1-score, confusion matric etc.

Dataset Description

- Implements a Support Vector Machine (SVM) linear and RBF classifiers to predict the class of breast tumor (breast cancer) from the provided dataset malignant or benign.
- The dataset needs to be loaded from scikit-learn
- The dataset should be used for training a SVM classifier and evaluate its performance using various metrics such as accuracy, precision, recall, F1-score

Dataset Details

• The breast cancer dataset consists of 569 samples, each representing a patient with a set of features

Data Source/Published By

- Breast Cancer Wisconsin (Diagnostic) UCI Machine Learning Repository
- Nuclear feature extraction for breast tumor diagnosis | Semantic Scholar
- Supported by CS Department, University of Wisconsin-Madison

Features & Data Download Link

Breast Cancer Wisconsin (Diagnostic) - UCI Machine Learning Repository

Project Implementation Demo







Interested in building a Gen Al application? Reach out to venkat@brillium.in

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