

## Different Types of SVM model-Contd.

- A. Kernel SVM:** Kernel SVM works on the dual of the SVM model such that it is a Similarity check between two data points  $x_1$  and  $x_2$  :

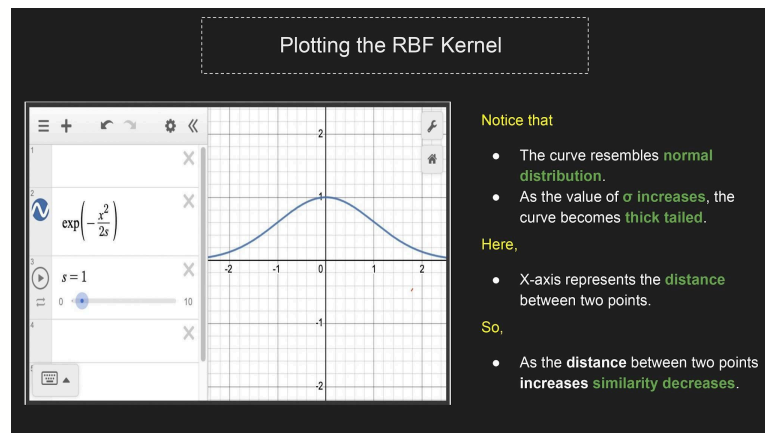
*Kernel Function* =  $k(x_1, x_2) = (x_1^t x_2 + c)^m$ , where m is the degree of Polynomial

### Understanding Kernel Trick

- **Kernel Trick:** Kernel function projects a d-dim data to d'- dim data (where  $d' \gg d$ ) such that the data points are easily separable
- Can use kernelization with LogReg and Deep Learning models also.

### Popular kernels

- Quadratic kernel: here  $m=2$ 
  - $K_q(x_1, x_2) = (x_1^t x_2 + c)^2$
  - Now if  $C=1$ , then on expanding  $K_q(x_1, x_2)$ , it becomes a 6 dimensional data.
- RBF(Radial Basis Function) / Gaussian kernel
  - Effective when not sure which degree of polynomial (m) to use
  - $K_{rbf}(x_1, x_2) = e^{\frac{-||x_1 - x_2||^2}{2\sigma^2}}$ ; where  $\sigma$  is a hyperparameter and  $||x_1 - x_2||^2$  is euclidean distance



### RBF kernel SVM vs kNN

- RBF kernel SVM is similar to KNN geometrically but differs in runtime complexity:  $O(\#SV * d)$ , where
  - #SV: Number of support vectors;
  - d: number of dimensions

## How is the Loss Function calculated for SVM?

The Loss function consists of two components:

- **Hinge Loss:** To have a minimum  $\zeta$ , such that  $y (w^T x + b) \geq 1 - \zeta$

$$\text{Hinge Loss: } \frac{1}{n} \sum_{i=1}^n \zeta_i$$

- **Max Margin:** To have a maximum Margin for generalizing on the data

$$\text{Max Margin: } \frac{\|w\|}{2}$$

Thus the total loss becomes:

$$\text{Loss: } \min_{(w,b)} \frac{\|w\|}{2} + C \frac{1}{n} \sum_{i=1}^n \zeta_i$$

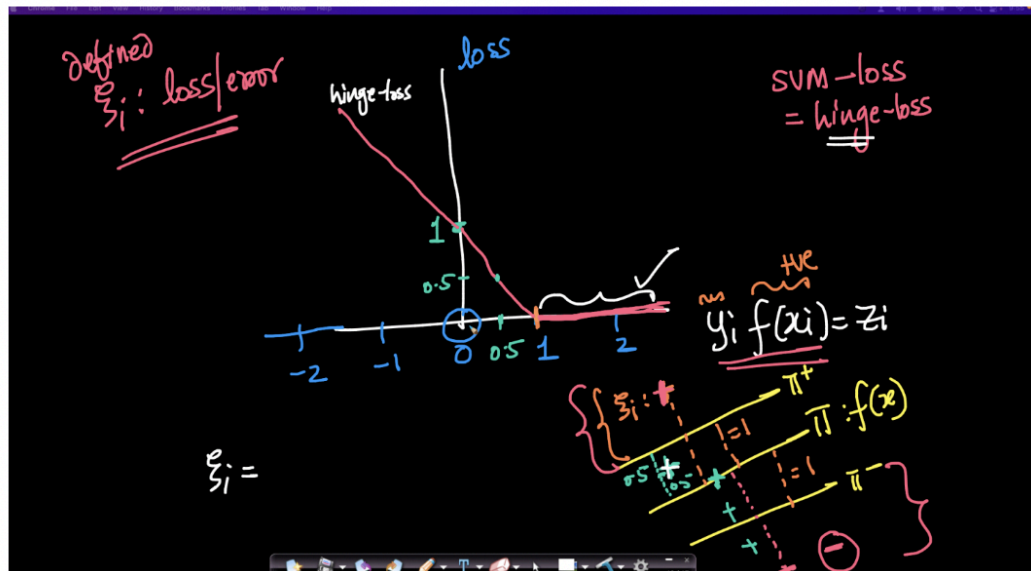
Where

- C is a hyperparameter that is analogous to the Regularization parameter ( $\lambda$ ) and -
- $\frac{\|w\|}{2}$  is analogous to L2 Regularization.

**Note:-**

- This is the **Primal form** of SVM
- The primal form works similarly to Log Reg.

## Understanding Hinge Loss



## Dual form of loss in SVM

We define a variable  $\alpha_i$  for each data point  $x_i$ , such that

- $0 \leq \alpha_i \leq C$
- $\sum_{i=1}^n \alpha_i y_i = 0$

Dual form of loss function:  $\max_{\alpha_i} \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_j y_i y_j x_i^T x_j$

**Note:** Dual form aids in the kernel trick by implicitly transforming the data points to a higher dimension

**Prediction using dual form:**

How do we predict the data points using dual form?

Prediction equation for dual form :-

$$\hat{y} = f(x_q) = \sum_{i=1}^N \alpha_i y_i x_i^T \cdot x_q$$

Diagram annotations:

- $\hat{y}$ : Query point
- $f(x_q)$ : Query point
- $\alpha_i$ : Learnt
- $y_i$ : Class label (-1 or +1)
- $x_i^T$ : Training data
- $x_q$ : Query point

Say, we have 10k data points.

## How does the value of C affect the model?

C is the hyperparameter. It causes a tradeoff between maximizing the margin and minimizing  $\zeta$

- If C is very large,
  - The SVM model tries to minimize HingeLoss
  - This makes the model have 0 incorrect predictions.
  - Thus making the model **overfit**.
- If C is very small,
  - The model tends to generalize the data
  - Hence, the model tends to have a maximum  $\zeta$ ,
  - Thus making the model **underfit**.

## What is the impact of Imbalanced Data on SVM?

SVM is impacted if there is an imbalance in Support Vectors. To resolve this:

- Either class weights should be used.
- Or rebalance the data.

## What are some Limitations of the SVM model?

- In some situations, the RBF kernel is very similar to KNN.
- In practice, GBDT/Random Forest still beats SVM
- Time complexity to train SVM is very high:  $O(n^2)$ ; n: no of data points

- Unlike Deep learning, SVM cannot create new features on its own.
- If we observe, we are just replacing feature engineering in GBDT/Random forest with kernel design in SVM.
- Even the RBF kernel SVM is impacted by outliers.

## Support Vector Regressor (SVR)

Though not very popularly used, SVM can be used for regression problems also.

Loss function:  $\min_{(w,b)} \frac{1}{2} ||w||^2 + C \cdot \epsilon$ , such that:

- $y_i - \hat{y}_i \leq \epsilon$
- $\hat{y}_i - y_i \leq \epsilon$
- $\epsilon \geq 0$