Applying Machine Learning to Image Data

Detroit Data Science
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Exercise: Image Orientation Detection

 0°
 90°
 180°
 270°

 180°
 180°
 180°
 180°
 180°

Exercise: Image Orientation Detection

- 4 possible classes: 0°, 90°, 180°, 270°
- Image features will be provided for you
 - Histogram of Oriented Gradients
 - Spatial color moments (3 mean and 3 variance values of L, U, and V)
 - Normalized spatial color moments
 - Principal Component Analysis
 - Linear Discriminant Analysis
- // TODO: Choose the combination of features and parameter values for training an SVM for classification

Training Data: 2149 images of people on bikes

























Testing Data: 626 images of people running























Exercise: What's included?

- Python code: function_list.py and student.py
- Random assignment and rotation of training and test images with extracted features:
 - /img_idx/trainingdata.pckl and /img_idx/testingdata.pckl image indices
 - /features/trainingdata_*_8x8.pckl training data features
 - /features/testingdata_*_8x8.pckl testing data features
- Conference and journal papers which concern image orientation detection: /papers/*.pdf
- An overview SVMs & the feature extraction approaches: /ppt/features.pdf

Overview: Histogram of Oriented Gradients

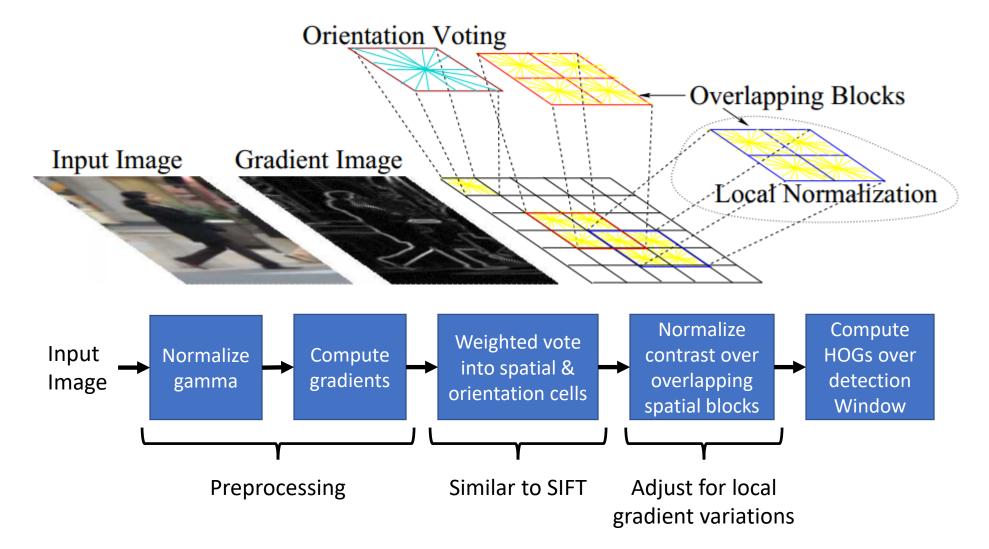


Image from Dalal and Triggs, a good tutorial can be found here: http://www.learnopencv.com/histogram-of-oriented-gradients/

HOG Preprocessing

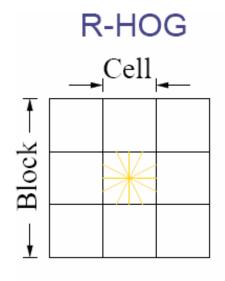
- Normalize over gamma for luminance correction
- Calculate gradient with mask [-1, 0, 1]
 - Larger masks (e.g., Sobel or 2-d Gaussian filtering) perform worse
 - Smoothing damages performance significantly
 - Color images have gradients per channel where the channel with the largest norm is used

HOG Spatial and Orientation Cells

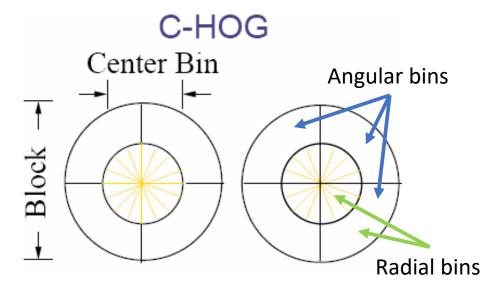
 Each pixel calculates a weighted vote based on the orientation of the gradient centered on it

- Calculate pixel orientation and magnitude
 - Orientation binned every 10° for 9 bins between $0^\circ \to 180^\circ$ (unsigned) or 18 bins between $0^\circ \to 360^\circ$ (signed)
 - Magnitude of the gradient = weight

HOG Contrast Normalization



 3×3 cell with 6×6 pixels per cell works well, though depends on the size of the object being detected



Good parameters are 4 angular bins (adding more worsens results) with 2 radial bins (more is unnecessary), and a center bin has a radius of 4 pixels

Normalization

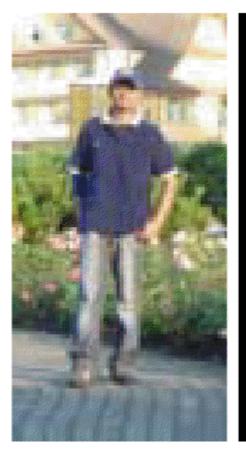
$$L1 - norm : v \longrightarrow v/(||v||_1 + \epsilon)$$

$$L2-norm: v \longrightarrow v/\sqrt{||v||_2^2+\epsilon^2}$$

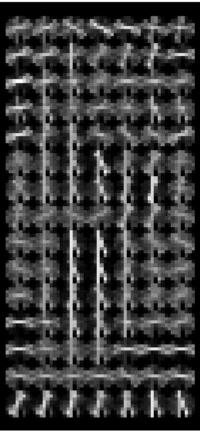
$$L1 - sqrt : v \longrightarrow \sqrt{v/(||v||_1 + \epsilon)}$$

L2-hys: L2-norm, plus clipping at .2 and renomalizing

HOG Example



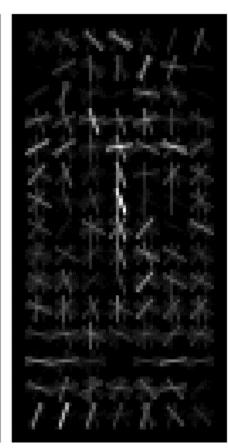
Image



R-HOG descriptor



Positive SVM weights



Negative SVM weights

HOG Examples (training set)

HOG feature extraction procedure:

- Resized to 200×200 pixels and converted to from RGB to grayscale
- Gaussian blur applied ($\sigma = 2$)
- HOG computation parameters:
 - Rectangular cell size of 20×20 pixels (for a 10×10 cell configuration)
 - L1 normalization over a 1×1 cell block
 - 4 orientations are used for binning

Original



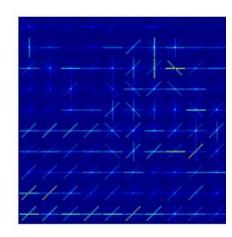


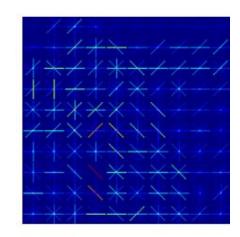
Adjusted





HOG

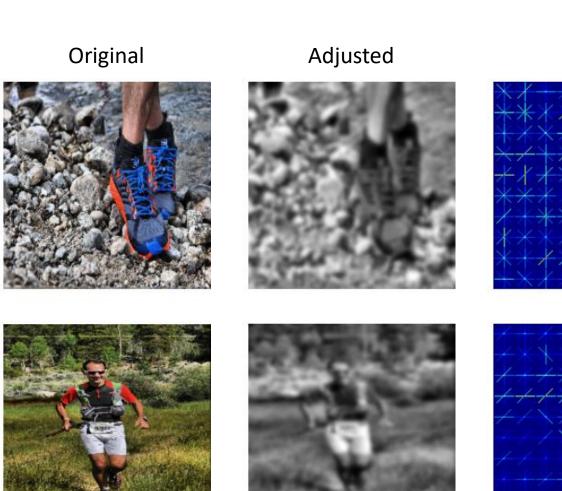




HOG Examples (test set)

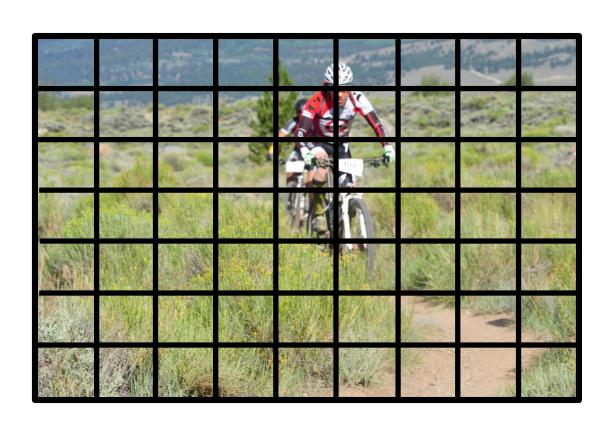
HOG feature extraction procedure:

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HOG

Overview: Spatial Color Moments



- Divide into blocks
- Compute mean and variance of the pixel values within each block for each color channel
- Concatenate all mean and variance values into a vector

Overview: Normalized Spatial Color Moments

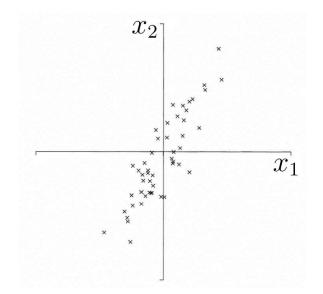
• Normalize the spatial color moment features to the same scale:

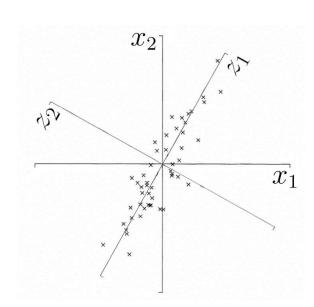
$$y_i' = \frac{y_i - min_i}{max_i - min_i}$$

- y_i = the i-th feature component of a feature vector y
- min_i , max_i = the range of values for the i-th feature component over the training samples
- All values are computed using L, U, V color space

Overview: Principal Component Analysis

- Multivariate statistical procedure explaining covariance structure of a set of variables via small number of their linear combinations.
- Determine axis with greatest variation for a given data set.
- Final coordinate system best represents the variance.





PCA

- Algebra Principal components are particular linear combinations of original variables, forming the projection that best represents the data with low mean square error (MSE)
- Geometry Linear combinations represent new coordinate system (via translations and rotations). New axes can determine directions of max variability of data set.
- Computational Principal components found by calculating eigenvectors and eigenvalues of data covariance matrix. Eigenvector with largest eigenvalue reveals the direction of greatest variation.

PCA

Formulate the mean adjusted data matrix.

$$U = \begin{pmatrix} x_{1,1} - \mu_1 & \cdots & x_{1,M} - \mu_M \\ \vdots & \ddots & \vdots \\ x_{N,1} - \mu_1 & \cdots & x_{N,M} - \mu_M \end{pmatrix}$$

Obtain the Covariance matrix.

$$\sum = U^T U / (N - 1)$$

Compute the projection matrix, each column being a direction of the new axes (an eigenvector).

$$\Phi = [\varphi_1, \varphi_2, \varphi_3, ..., \varphi_L]$$
 where $L \leq M$

PCA

• The set of L eigenvectors of the covariance corresponds to the L largest eigenvalues that minimizes the MSE of reconstruction over all choices of an orthonormal basis of size L.

- Much of the variability of data set can be accounted for in a smaller number of L principal components (pcs) as opposed to potentially large M variables.
- N examples of M variables has suddenly been reduced to N examples on L pcs.

PCA Example



PCA Example

• Our Mean Image:



• Eigenvectors (also known as eigenfaces):









PCA Example

• Reconstruct an Image from eigen-basis











PCA Summary

Use eigenvectors of covariance matrix with largest eigenvalues – projects across dimensions of maximum variance

• Is a linear transform of high dimensional data to a lower dimension whose components are uncorrelated (if data is Gaussian, uncorrelated implies statistical independence)

• Is optimal when the you want to minimize the approximated mean square error of the projection

PCA Summary

 Works well if the data points are distributed throughout the hyperplane

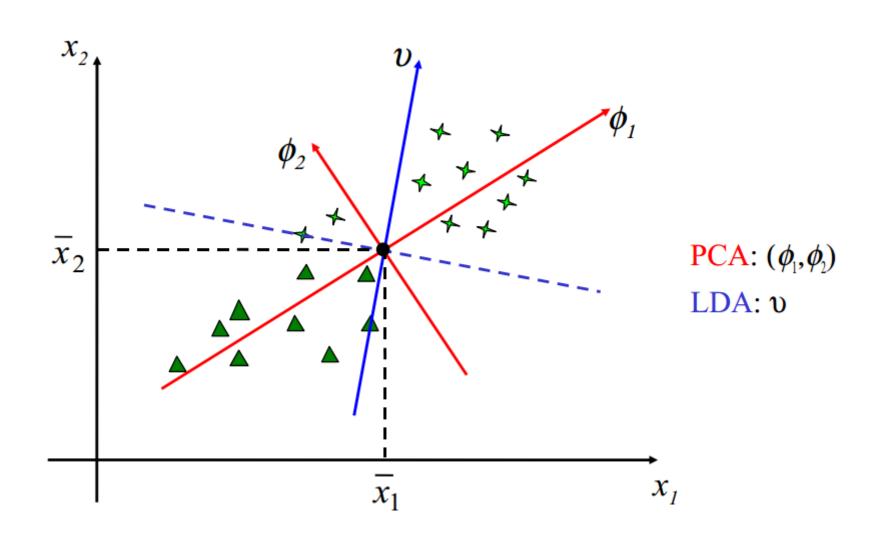
 Does not necessarily lead to good class separability – no guarantee that the components will be discriminable features

 Will fail if data is highly non-linear or lies on more complicated manifolds

Overview: Linear Discriminant Analysis

- Statistical pattern recognition technique.
- Separates data into discrete groups via transformation into a space that maximizes 'between-class' separation while minimizing their 'with-in class' variability.
- Looks for a projection where new directions maximally separate data.

LDA Geometric View



LDA

- Within Class scatter matrix $\rightarrow S_w = \sum_{i=1}^c \sum_{j=1}^{n_i} (Y_j \mu_i) (Y_j \mu_i)^T$
- Between Class scatter matrix $\rightarrow S_b = \sum_{i=1}^{c} (\mu_i \mu) (\mu_i \mu)^T$
- The goal for LDA is to find a projection matrix that maximizes the ratio of determinant of S_b to the determinant of S_w (Fisher's Criterion)

$$\mathbf{\Phi}_{LDA} = \max |\mathbf{\Phi}^T S_b \mathbf{\Phi}| / |\mathbf{\Phi}^T S_w \mathbf{\Phi}|$$

• If S_w is non-singular then Fisher's Criterion is maximized when the projection matrix comprises of eigenvectors from $S_w^{-1}S_b$

LDA Summary

• LDA is a statistical pattern recognition technique for separating samples into discrete groups.

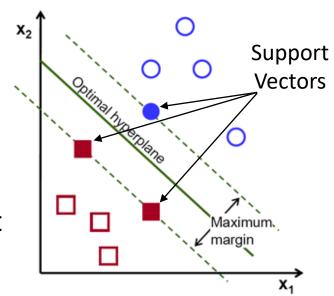
• LDA looks for directions that are efficient for *discriminating* data whereas PCA looks for data that are efficient for *representing* data.

• When the training data set is small, PCA can outperform LDA, also PCA is less sensitive to a different training data set.

- Designed to find the maximum margin between two classes in feature space
- The optimal hyperplane (decision surface) is one that separates our data with the largest margin between the features of each class
- 'Support Vectors' are the data points that lie closest to the hyperplane

Weight vector
$$f(x) = \psi^T x + \beta \begin{cases} \geq 0 & \text{Class A} \\ < 0 & \text{Class B} \end{cases}$$
Input vector Bias

• Classifying a feature is simply labeling whether the new point falls on one side or the other of the trained margin



- w is orthogonal to x when x is located on the hyperplane (i.e., $w^T x_{\perp} + \beta = 0$)
- By subtracting the two planes we can find our margin (separation between points x_m and x_p):

$$w^{T}x_{p} + \beta = 1$$

$$-w^{T}x_{m} + \beta = -1$$

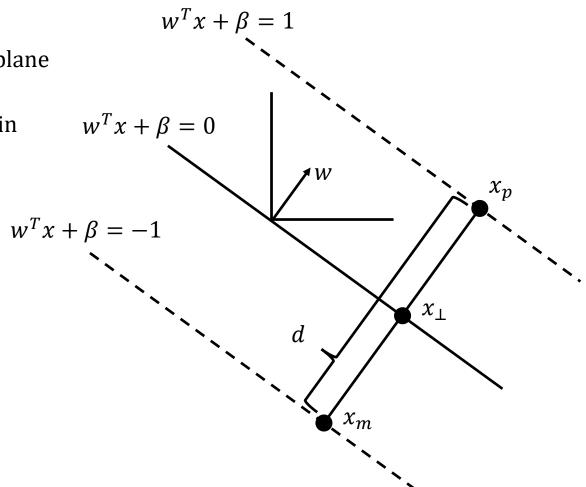
$$w^{T}(x_{p} - x_{m}) = 2$$

Since w is perpendicular, the closest point to x_m can be defined as:

$$x_p = x_m + dw$$

$$\therefore dw^T w = 2 \to d = \frac{2}{w^T w} = \frac{2}{\|w\|_2^2}$$

• Distance = $d||w||_2 = \frac{2}{||w||_2^2} ||w||_2 = \frac{2}{\sqrt{w^T w}}$



 $dw = \text{line segment connecting } x_m \text{ and } x_p$ $d||w||_2 = \text{distance between } x_m \text{ and } x_p$

- We want to maximize the margin, i.e. the distance between two class boundaries $\frac{2}{\sqrt{w^T w}}$, which is equivalent to minimizing $\frac{\sqrt{w^T w}}{2}$ (which is equivalent to minimizing $\frac{1}{2} w^T w$ since the square root function is monotonic)
- Additionally, we can define our labels $y_i = 1 \ \forall i \in \text{Class A}$ and $y_i = -1 \ \forall i \in \text{Class B}$:

Class A:
$$w^T x_i + \beta \ge 1$$
, $y_i = 1 \to y_i (w^T x_i + \beta) \ge 1$
Class B: $w^T x_i + \beta \le -1$, $y_i = -1 \to y_i (w^T x_i + \beta) \ge 1$

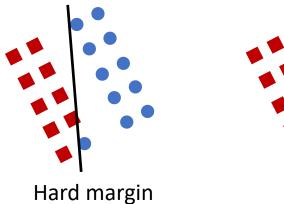
Accordingly, the convex optimization problem for a linear SVM with a hard margin:

$$\min_{w,\beta} \frac{1}{2} w^T w \text{ subject to: } y_i(w^T x_i + \beta) \ge 1 \ \forall i$$

- In practice, training samples may contain noise and are not linearly separable, giving no feasible solution to the hard margin problem
- The soft margin SVM uses a 'slack variable', ε_i , to represent the error of the i-th training sample (constrained optimization problem):

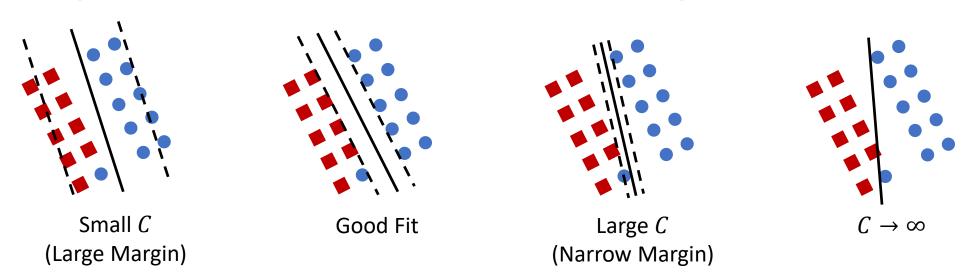
$$\min_{w,\beta,\varepsilon} w^T w + 2C \sum_{i} \varepsilon_i \text{ subject to: } y_i(w^T x_i + \beta) \ge 1 - \varepsilon_i \text{ and } \varepsilon_i \ge 0 \ \forall i$$

A regularization parameter that weighs the cost of the penalty for misclassification. Determined using cross-validation ($C \rightarrow \infty$ = hard margin)

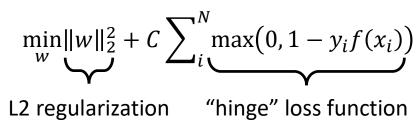


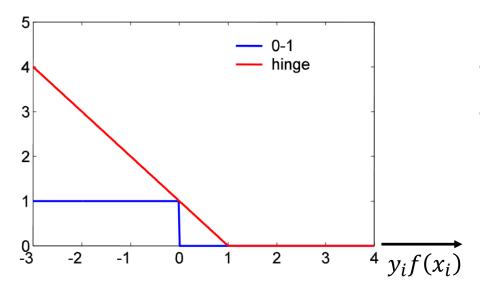


- ullet As a regularization parameter, C, determines the penalty for misclassifications
 - Large C makes constraints hard to ignore (narrow margin) and will lead to a lower bias, higher variance SVM (i.e., overfitting)
 - Small C allows for constraints to be ignored (large margin) and will lead to a higher bias, lower variance SVM (i.e., underfitting)



- Learning an SVM was previously formulated as a constrained optimization problem over w and ε , however, $y_i(w^Tx_i+\beta) \geq 1-\varepsilon_i = y_if(x_i) \geq 1-\varepsilon_i$ and with $\varepsilon_i \geq 0$, can be more concisely written as: $\varepsilon_i = \max(0, 1-y_if(x_i))$ Known as the 'Primal Problem'
- The equivalent learning problem is unconstrained over w:





- The function takes the form of arg min $\sum \mathcal{L}(y, f(x)) + \lambda R(f)$
- The graph represents two margin based classifiers which induce a decision rule via sign(f):
 - Misclassification (0-1) loss: $\mathcal{L}(y, f(x)) = I(yf(x) \le 0)$
 - Hinge loss (SVM): $\mathcal{L}(y, f(x)) = \max(0, 1 yf(x))$

- Problem: the primal problem is convex but not differentiable
- Solution: the 'Dual problem' can be derived using Lagrangian multipliers

$$\max_{\alpha} \sum_{i} \alpha_{i} - \frac{1}{2} \sum_{jk} \alpha_{j} \alpha_{k} y_{j} y_{k} (x_{j}^{T} x_{k}) \text{ subject to: } 0 \leq \alpha_{i} \leq C \text{ and } \sum_{i} \alpha_{i} y_{i} = 0$$

- The dual form benefits:
 - Requires to need to learn N parameters, while the primal form requires learning δ parameters ($x_i \in \mathbb{R}^{\delta}$ for $i=1\cdots N$, if $N\ll \delta$ then it's more efficient to solve for α than w)
 - Only involves $(x_i^T x_k)$ which is helpful for non-linear decision boundaries



• Given a non-linear decision boundary, the data may be linearly separable in a different space, whereby $\phi(x)$ is a feature map $(\phi: x \to \phi(x) \mathbb{R}^{\delta} \to \mathbb{R}^{D})$:

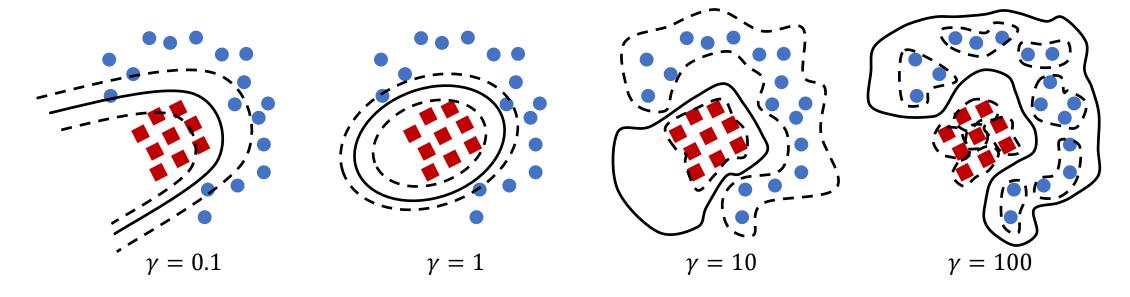
$$f(x) = w^T \phi(x) + \beta$$

- Map x to $\phi(x)$ where the data is separable (solving for w in the high dimensional space \mathbb{R}^D)
- If $D \gg \delta$? Use the dual formulation: $k(x_j, x_k) = \phi(x_j)^T \phi(x_k)$ where $k(x_j, x_k)$ is the 'Kernel'

Common Kernels:

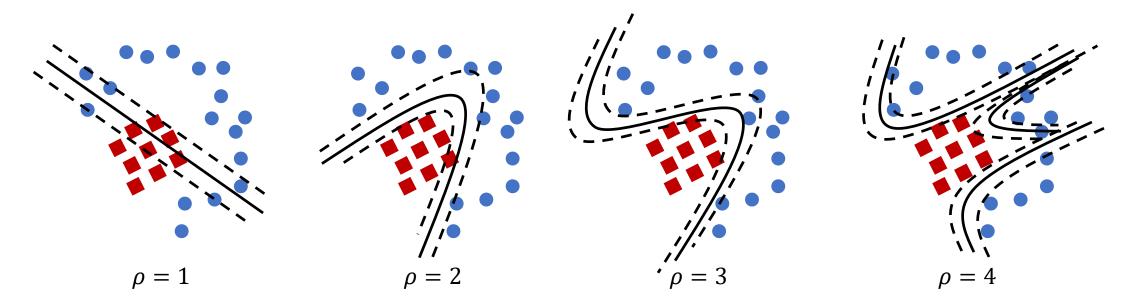
- <u>Linear kernel</u>: $k(x, y) = x^T y$ (no mapping)
- Polynomial kernel: $k(x,y) = (\gamma x^T y + r)^{\rho}$ for $\gamma, \rho > 0$ and $r \ge 0$ (polynomial terms up to degree ρ)
- Radial basis function (RBF) kernel (Gaussian): $k(x,y) = \exp\left(-\frac{\|x-y\|_2^2}{2\sigma^2}\right) = \exp(-\gamma \|x y\|_2^2)$ where $\gamma = \frac{1}{2\sigma^2}$, the RBF kernel is shift invariant $(k(x+a,y+a) = k(x,y) \ \forall a$
- Sigmoid kernel: $k(x, y) = \tanh(\gamma x^T y + r)$

RBF Kernel: $k(x, y) = \exp(-\gamma ||x - y||_2^2)$



 γ is known as the kernel bandwidth, as γ increases, the flexibility of the decision boundary increases. Small values produce a boundary that is near linear and as γ increases the SVM overfits.

Polynomial kernel: $k(x, y) = (\gamma x^T y + r)^{\rho}$ $\rho = \text{degree}, \gamma = \text{scale}, r = \text{bias (offset)}$



The polynomial degree, ρ , controls the flexibility of the decision boundary increases. When $\rho=1$ the boundary is linear, and as ρ increases the SVM will often overfit.

Grid Search

 As the number of hyperparameters increases, the complexity of choosing acceptable values also increases

Grid search iterates over varying values of each hyperparameter while

holding all others constant

 Grid points are usually chosen on a logarithmic scale and classifier accuracy is estimated for each point on the grid

Accuracy at $\gamma=10^{-3}$ and $C=10^{1}$

