Kernel Methods for Pattern Recognition

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• **Examples** \( (x_i) \)

\[ \text{fMRI/sMRI} \]
\[ \text{3D matrix of voxels} \]

\[ x_i \] is a vector of size \( dx1 \) where \( d \) is the number of voxels

• **Labels** \( (y_i) \)
  • categorical value for classification (e.g. class 1 = patients, class 2 = healthy controls)
  • continuous value for regression (e.g. age or clinical scale).

• **Matrix notation** (one example per row)

\[
X = [x_1 \ x_2 \ ... \ x_N]^T \\
y = [y_1 \ y_2 \ ... \ y_N]^T
\]
Pattern Recognition Framework

Learning/Training Phase
- Generate a function or classifier $f$ such that $f(x_i) \rightarrow y_i$

Testing Phase
- Prediction $f(x_i) = y_i$

Computer-based procedures that learn a function from a series of examples

Input (brain scans)
- $x_1$
- $x_2$
- $x_3$

Training Examples:
- $(x_1, y_1), \ldots, (x_s, y_s)$

Output (control/patient)
- $y_1$
- $y_2$
- $y_3$

No mathematical model available
Linear models

- Linear predictive models (classifier or regression) are parameterized by a weight vector $\mathbf{w}$ and a bias term $b$.
  
  \[ f(\mathbf{x}_*) = \mathbf{w} \cdot \mathbf{x}_* + b \]

  where $f(\mathbf{x}_*)$ is the **predicted score for regression** or the distance to the **decision boundary for classification models**.

- The weight vector can be expressed as a linear combination of training examples $\mathbf{x}_i$ (where $i = 1, \ldots, N$ and $N$ is the number of training examples).
  
  \[ \mathbf{w} = \sum_{i=1}^{N} \alpha_i \mathbf{x}_i \]
Main difficulties:
• Very high dimensional data: computational issues
• Often the dimensionality of the data is greater than the number of examples: ill-conditioned problems

Potential Solutions:
• Feature Selection
• Region of Interest
• Searchlight
• Kernel Methods + Regularisation -> PRoNTo
Kernel Methods

• The kernel methodology provides a powerful and unified framework for investigating general types of relationships in the data (e.g. classification, regression, etc).

• **Kernel methods consist of two parts:**
  ✓ Computation of the kernel matrix (mapping into the feature space).
  ✓ A learning algorithm based on the kernel matrix (designed to discover linear patterns in the feature space).

• **Advantages:**
  ✓ Represent a *computational shortcut* which makes possible to represent linear patterns efficiently in high dimensional space.
  ✓ Using the dual representation with proper *regularization* enables efficient solution of ill-conditioned problems.

* e.g. restricting the choice of functions to favor functions that have small norm.
Kernel Function ("similarity measure")

- Kernel is a function that, for given two pattern \( x \) and \( x^* \), returns a real number characterizing their similarity.

- A simple type of similarity measure between two vectors is a dot product (linear kernel).

\[
\text{Dot product} = (4\times-2)+(1\times3) = -5
\]

\[
K_{\text{linear}} = XX^T
\]
• There are more general “similarity measures”, i.e. nonlinear kernels: Gaussian kernel, Polynomial kernel, etc.

• Nonlinear kernels are used to map the data to a higher dimensional space as an attempt to make it linearly separable.

• The kernel trick enables the computation of similarities in the feature space without having to compute the mapping explicitly.
Advantage of linear models

- Neuroimaging data are extremely high-dimensional and the sample sizes are very small, therefore non-linear kernels often don’t bring any benefit.

- Linear models reduce the risk of overfitting the data and allow direct extraction of the weight vector as an image (i.e. predictive map).
Learning with kernels

- Making predictions with kernel methods

\[ f(x_*) = w \cdot x_* + b \]  \hspace{1cm} \text{Primal representation}

\[ f(x_*) = \sum_{i=1}^{N} \alpha_i x_i \cdot x_* + b \]

\[ f(x_*) = \sum_{i=1}^{N} \alpha_i K(x_i, x_*) + b \]  \hspace{1cm} \text{Dual representation}
How to interpret the weight vector ($w$)?

Examples of class 1

Examples of class 2

New example

Testing

Model weight vector

$\mathbf{w} = (w_1, w_2) = (+5, -3)$

Spatial representation of decision function

Multivariate pattern -> No local inferences should be made!

$\mathbf{f}(\mathbf{x}) = (w_1v_1 + w_2v_2) + b$

$= (+5 \cdot 0.5 - 3 \cdot 0.2) + 0$

$= 1.9$

Positive value -> Class 1
Examples of Kernel Methods in PRoNTo

- Support Vector Machines (SVM)
- Gaussian Processes (GP)
- Kernel Ridge Regression (KRR)
- Relevance Vector Regression (RVR)
- Multiple Kernel Learning (MKL)
Example of Kernel Methods

(1) Support Vector Machine

SVM became famous when, using images as input, it gave accuracy comparable to neural-network in a handwriting recognition task.

Currently, SVM is widely used in object detection & recognition, text recognition, biometrics, speech recognition, neuroimaging, etc.

Also used for regression.
• Among all hyperplanes separating the data there is a unique **optimal hyperplane**, the one which presents the **largest margin** (the distance of the closest points to the hyperplane).

• Let us consider that all test points are generated by adding bounded noise ($r$) to the training examples (test and training data are assumed to have been generated by the same underlying dependence).

• If the optimal hyperplane has margin $\rho > r$ it will correctly separate the test points.
• We assume that the data are linearly separable, that is, there exist \( \mathbf{w} \in \mathbb{R}^d \) and \( b \in \mathbb{R} \) such that \( y_i (\mathbf{w} \cdot \mathbf{x}_i + b) > 0 \), \( i = 1, \ldots, m \).

• Rescaling \( \mathbf{w} \) and \( b \) such that the points closest to the hyperplane satisfy \( |(\mathbf{w} \cdot \mathbf{x}_i + b)| = 1 \) we obtain the canonical form of the hyperplane satisfying \( y_i (\mathbf{w} \cdot \mathbf{x}_i + b) > 0 \)

• The distance of a point \( \mathbf{x}_i \) to a hyperplane \( H_{\mathbf{w}, b} \) is given by \( \rho_x = |(\mathbf{w} \cdot \mathbf{x}_i + b)|/||\mathbf{w}|| \)

• The distance from the closest point to the canonical hyperplane is \( \rho = 1/||\mathbf{w}|| \).

• In this case, the margin, measured perpendicularly to the hyperplane, equals \( 2/||\mathbf{w}|| \).

• In order to maximize the margin we need to minimize \( ||\mathbf{w}||/2 \).
Linearly separable case (Hard Margin SVM)

• Constrained optimization problem

\[
\begin{align*}
\min & \quad \frac{1}{2} \| w \|^2 \\
\text{s.t.} & \quad y_i(w \cdot x_i + b) \geq 1, \ i = 1, \ldots, m
\end{align*}
\]

Quadratic problem: unique optimal solution

• The solution of this problem is equivalent to determine the saddle point of the Lagrangian function

\[
L(w, b; \alpha) = \frac{1}{2} \| w \|^2 - \sum_{i=1}^{N} \alpha_i \left\{ y_i (w \cdot x_i + b) - 1 \right\}
\]

where \( \alpha_i \geq 0 \) are the Lagrange multipliers.

• We minimize \( L \) over \((w,b)\) and maximize over \( \alpha \).
Differentiating $L$ w.r.t. $w$ and $b$ we obtain:
\[
\frac{\partial L}{\partial b} = - \sum_{i=1}^{N} y_i \alpha_i = 0
\]
\[
\frac{\partial L}{\partial w} = w - \sum_{i=1}^{N} \alpha_i y_i x_i = 0 \implies w = \sum_{i=1}^{N} \alpha_i y_i x_i
\]

Substituting $w$ in $L$ leads to the dual problem
\[
\max \ Q(\alpha) := -\frac{1}{2} \alpha^T A \alpha + \sum_{i=1}^{N} \alpha_i
\]
\[
\text{s.t. } \sum_{i} y_i \alpha_i = 0
\]
\[
\alpha_i \geq 0, \ i=1,\ldots,N
\]

where $A$ is an $N \times N$ matrix and $A = (y_i y_j x_i \cdot x_j : i, j = 1,\ldots,N)$.
If $\alpha$ is a solution of the dual problem then the solution $(w, b)$ of the primal problem is given by

$$w = \sum_{i=1}^{N} \alpha_i y_i x_i$$

Note that $w$ is a linear combination of only the $x_i$ for which $\alpha_i > 0$. These $x_i$ are called support vectors (SVs).

Parameter $b$ can be determined by $b = y_i - w \cdot x_i$, where $x_i$ corresponds to a SV.

A new point $x_\star$ is classified as

$$f(x_\star) = \text{sgn}\left(\sum_{i=1}^{N} y_i \alpha_i x_i \cdot x_\star + b\right)$$

The dot product is simple type of similarity measure.
Kernel Trick

- The dot product can be replaced by a kernel function which corresponds to a dot product in the feature space.

- The kernel trick is a way of mapping observations from the original space into a feature space, without ever having to compute the mapping explicitly.
Some remarks

• The fact that the Optimal Separating Hyperplane is determined only by the support vectors is most remarkable. Usually, the support vectors are a small subset of the training data.

• All the information contained in the data set is summarized by the support vectors. The whole data set could be replaced by only these points and the same hyperplane would be found.
• If the data is not linearly separable the previous analysis can be generalized by looking at the problem

$$\min \frac{1}{2} \| w \|^{2} + C \sum_{i=1}^{N} \xi_{i}$$

s.t. $y_{i}(w \cdot x_{i} + b) \geq 1 - \xi_{i}$

$$\xi_{i} \geq 0, \quad i = 1, ..., N$$

• The idea is to introduce the slack variables $\xi_{i}$ to relax the separation constraints ($\xi_{i} > 0 \Rightarrow x_{i}$ has margin less than 1).
New dual problem

• A saddle point analysis (similar to that above) leads to the dual problem

\[ \max Q(\alpha) := -\frac{1}{2} \alpha^T A \alpha + \sum_i \alpha_i \]

\[ \text{s.t. } \sum_t y_i \alpha_i = 0 \]

\[ 0 \leq \alpha_i \leq C, \quad i = 1, \ldots, N \]

• This is like the previous dual problem except that now we have “box constraints” on \( \alpha_i \).

• Again we have \( \mathbf{w} = \sum_{i=1}^{N} \alpha_i y_i \mathbf{x}_i \)
The role of the parameter C

• The parameter C that controls the relative importance of minimizing the norm of $\mathbf{w}$ (which is equivalent to maximizing the margin) and satisfying the margin constraint for each data point.

• If C is close to 0, then we don't pay that much for points violating the margin constraint. This is equivalent to creating a very wide tube or safety margin around the decision boundary (but having many points violate this safety margin).

• If C is close to $\infty$, then we pay a lot for points that violate the margin constraint, and we are close the hard-margin formulation we previously described - the difficulty here is that we may be sensitive to outlier points in the training data.

• C is often selected by cross-validation (nested cross-validation in PRoNTo).
• SVMs are prediction devices known to have good performance in high-dimensional settings.

• "The key features of SVMs are the use of kernels, the absence of local minima and the sparseness of the solution.” Shawe-Taylor and Cristianini (2004).
Example of Kernel Methods

(1) Multiple Kernel Learning (MKL)
• Many practical learning problems involve multiple, heterogeneous data sources.

• It seems advantageous to combine different sources of information for prediction (e.g. multimodal imaging for diagnosis/prognosis).

• Need to learn with not only a single kernel but with multiple kernels.
Multiple Kernel Learning (MKL) has been proposed as an approach to simultaneously learn the kernel weights and the associated decision function in supervised learning settings.

In MKL, the kernel $K$ can be considered as a linear combination of $M$ “basis kernels”

$$K(x, x') = \sum_{i=1}^{M} d_m K_m(x, x')$$

with $d_m \geq 0$, $\sum_{i=1}^{M} d_m = 1$

The decision function of an MKL problem can be then expressed in the form:

$$f(x_s) = \sum_{i=1}^{m} w_m \cdot x_s + b$$
• One example of MKL approach based on SVM is the SimpleMKL (Rakotomamonjy, et al. 2008).

• SimpleMKL optimization problem

\[
\begin{align*}
\min & \frac{1}{2} \sum_{m=1}^{M} \frac{1}{d_m} \| w_m \|^2 + C \sum_{i=1}^{N} \xi_i \\
\text{s.t.} & \quad y_i \left( \sum_{m=1}^{M} w_m \cdot x_i + b \right) \geq 1 - \xi_i, \quad i = 1, \ldots, N \\
\xi_i & \geq 0, \quad i = 1, \ldots, N \\
\sum_{m=1}^{M} d_m & = 1, \quad d_m \geq 0, \quad m = 1, \ldots, M
\end{align*}
\]

the L1 constrain on \( d_m \) enforces sparsity on the kernels with a contribution to the model.
Single vs. Multiple Kernel Learning

Single kernel SVM

\[ f(x_*) = w \cdot x_* + b \]

Multiple kernel SVM

\[ f(x_*) = \sum_{i=1}^{m} d_i f_i(x_*) \]

\[ = \sum_{i=1}^{m} d_i w_i \cdot x_* + b \]

where \( d_i \) are weights and \( f_i(x_*) \) are individual kernel functions.
Example of Kernel Methods

(3) Kernel Ridge Regression
Kernel Methods

- The general equation for making predictions with kernel methods is

\[ f(x_*) = w \cdot x_* + b = \sum_{i=1}^{N} \alpha_i x_i \cdot x_* + b = \sum_{i=1}^{N} \alpha_i K(x_i, x_*) + b \]

where \( f(x_*) \) is the predicted score for regression or the distance to the decision boundary for classification.

- \( \alpha_i \) is the dual weight vector and \( b \) is a constant offset, both of which are learnt from the training samples.

- We can simplify the equation for making predictions by adding a constant element to \( x_* \), so that \( x_* = [x_* \: 1]^T \) and \( w=[w \: b]^T \)

\[ f(x_*) = w \cdot x_* \]
Kernel Ridge Regression

- Kernel ridge regression is the dual representation of ridge regression, which is sometimes known as the linear Least Square Regression (LSR) with Tikhonov regularization (Chu et al. 2011).

Illustration of a linear least squares fitting with $X \in \mathbb{R}^2$. We seek the linear function of $X$ that minimizes the sum of squared residuals from $Y$.

Hastie, Tibshirani & Friedman, 2009
Least Squares Regression (LSR)

- In LSR we compute the weight vector $\mathbf{w}$ by minimizing the mean squared errors on all training examples:

$$
\mathbf{w}^* = \text{argmin}_w \frac{1}{N} \sum_{i=1}^{N} (x_i \cdot \mathbf{w} - y_i)^2
$$

Using a matrix notation where $\mathbf{X} = [x_1 \ x_2 \ldots \ x_N]^T$ is a matrix containing the training examples vectors as its row we can rewrite the cost function as

$$
\mathbf{w}^* = \text{argmin}_w (\mathbf{Xw} - \mathbf{y})^T (\mathbf{Xw} - \mathbf{y})
$$

- To find the optimum $\mathbf{w}$ we set the derivative of the cost function with respect to $\mathbf{w}$ to 0, which yields to the following equation:

$$
\mathbf{X}^T (\mathbf{Xw} - \mathbf{y}) = 0
$$

$$
\mathbf{X}^T \mathbf{X} \mathbf{w} = \mathbf{X}^T \mathbf{y}
$$

$$
\mathbf{w} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}
$$
When the sample size is limited, i.e. in order to solve ill-posed problems or to prevent over-fitting, some form of regularization is often introduced into the model.

The regularization parameter $\lambda$ controls the amount of regularization.

Setting the derivative of the cost function with respect to $w$ to 0, which yields to the following equation:

$$w^* = \arg\min_w \frac{1}{N} \sum_{i=1}^{N} (x_i \cdot w - y_i)^2 + \lambda \|w\|^2$$

$$X^T (Xw - y) + \lambda w = 0$$

$$(X^T X + \lambda I)w = X^T y$$

$$w = (X^T X + \lambda I)^{-1} X^T y$$
Consider the general equation for making predictions

\[ f(x_*) = w \cdot x_* \]

\[ w = \sum_{i=1}^{N} \alpha_i x_i \]

To estimate the weights \( w \) we seek to minimize the empirical risk which is penalized to restrict model flexibility

\[ w^* = \arg\min_w \frac{1}{N} \sum_{i=1}^{N} L(y_i, x_i, w) + \lambda J(w) \]
Statistical Learning – General Framework

\[ w^* = \arg\min_w \frac{1}{N} \sum_{i=1}^{N} L(y_i, x_i, w) + \lambda J(w) \]

- **Loss function**: denotes the price we pay when we make mistakes in the predictions (e.g. squared loss, Hinge loss).

- **Regularization term**: favours certain properties and improves the generalisation over unseen examples (e.g. L2-norm, L1-norm).

- Many learning algorithms are particular choices of \( L \) and \( J \) (e.g. SVM, Kernel Ridge Regression).

\[
\text{KRR} \quad w^* = \arg\min_w \frac{1}{N} \sum_{i=1}^{N} (x_i \cdot w - y_i)^2 + \lambda \|w\|^2
\]

\[
\text{SVM} \quad w^* = \arg\min_w C \frac{1}{N} \sum_{i=1}^{N} \max[1 - y_i (x_i \cdot w + b), 0] + \lambda \|w\|^2
\]
<table>
<thead>
<tr>
<th>Model</th>
<th>Accuracy (%)</th>
</tr>
</thead>
<tbody>
<tr>
<td>LASSO</td>
<td>86.31%</td>
</tr>
<tr>
<td>Elastic Net</td>
<td>88.02%</td>
</tr>
<tr>
<td>Total Variation (TV)</td>
<td>85.79%</td>
</tr>
<tr>
<td>Laplacian (LAP)</td>
<td>83.71%</td>
</tr>
<tr>
<td>Sparse TV</td>
<td>85.86%</td>
</tr>
<tr>
<td>Sparse LAP</td>
<td>87.05%</td>
</tr>
</tbody>
</table>

- Weight maps for classifying fMRI images during visualization of pleasant vs. unpleasant pictures.
- All models used a square loss + regularization.

References


