

Kernel Methods

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Big Feature Spaces for Linear Models

The Input Space \mathcal{X}

- Our general learning theory setup: no assumptions about \mathcal{X}
- But $\mathcal{X} = \mathbf{R}^d$ for the specific methods we've developed:
 - Ridge regression
 - Lasso regression
 - Support Vector Machines
- Our hypothesis space for these was all affine functions on \mathbf{R}^d :

$$\mathcal{F} = \{x \mapsto w^T x + b \mid w \in \mathbf{R}^d, b \in \mathbf{R}\}.$$

- What if we want to do prediction on inputs not natively in \mathbf{R}^d ?

The Input Space \mathcal{X}

- Often want to use inputs not natively in \mathbf{R}^d :
 - Text documents
 - Image files
 - Sound recordings
 - DNA sequences
- But everything in a computer is a sequence of numbers?
 - The i th entry of each sequence should have the same “meaning”
 - All the sequences should have the same length

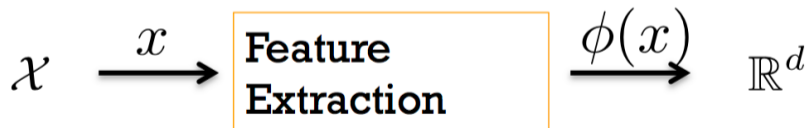
Feature Extraction

Definition

Mapping an input from \mathcal{X} to a vector in \mathbb{R}^d is called **feature extraction** or **featurization**.

Raw Input

Feature Vector

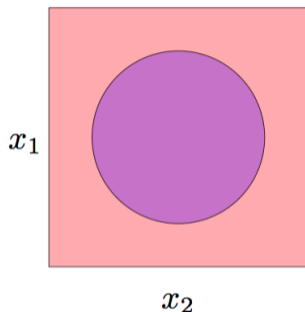


Linear Models with Explicit Feature Map

- Input space: \mathcal{X} (no assumptions)
- Introduce **feature map** $\psi : \mathcal{X} \rightarrow \mathbf{R}^d$
- The feature map maps into the **feature space** \mathbf{R}^d .
- Hypothesis space of affine functions on feature space:

$$\mathcal{F} = \{x \mapsto w^T \psi(x) + b \mid w \in \mathbf{R}^d, b \in \mathbf{R}\}.$$

Geometric Example: Two class problem, nonlinear boundary



- With identity feature map $\psi(x) = (x_1, x_2)$ and linear models, can't separate regions
- With appropriate featurization $\psi(x) = (x_1, x_2, x_1^2 + x_2^2)$, becomes linearly separable .
- Video: <http://youtu.be/3liCbRZPrZA>

From Percy Liang's "Lecture 3" slides from Stanford's CS221, Autumn 2014.

Expressivity of Hypothesis Space

- For linear models, to grow the hypothesis spaces, we must add features.
- Sometimes we say a larger hypothesis is “more expressive”.
 - (can fit more relationships between input and action)
- Many ways to create new features.

Handling with Nonlinearity with Linear Methods

Example Task: Predicting Health

- General Philosophy: Extract every feature that might be relevant
- Features for medical diagnosis
 - height
 - weight
 - body temperature
 - blood pressure
 - etc...

From Percy Liang's "Lecture 3" slides from Stanford's CS221, Autumn 2014.

Feature Issues for Linear Predictors

- For linear predictors, it's important **how** features are added
- Three types of nonlinearities can cause problems:
 - ① Non-monotonicity
 - ② Saturation
 - ③ Interactions between features

From Percy Liang's "Lecture 3" slides from Stanford's CS221, Autumn 2014.

Non-monotonicity: The Issue

- Feature Map: $\phi(x) = [1, \text{temperature}(x)]$
- Action: Predict health score $y \in \mathbf{R}$ (positive is good)
- Hypothesis Space $\mathcal{F} = \{\text{affine functions of temperature}\}$
- Issue:
 - Health is not an affine function of temperature.
 - Affine function can either say
 - Very high is bad and very low is good, or
 - Very low is bad and very high is good,
 - But here, both extremes are bad.

From Percy Liang's "Lecture 3" slides from Stanford's CS221, Autumn 2014.

Non-monotonicity: Solution 1

- Transform the input:

$$\phi(x) = \left[1, \{\text{temperature}(x) - 37\}^2 \right],$$

where 37 is “normal” temperature in Celsius.

- Ok, but requires manually-specified domain knowledge
 - Do we really need that?

Non-monotonicity: Solution 2

- Think less, put in more:

$$\phi(x) = \left[1, \text{temperature}(x), \{\text{temperature}(x)\}^2 \right].$$

- **More expressive** than Solution 1.

General Rule

Features should be simple building blocks that can be pieced together.

Saturation: The Issue

- Setting: Find products relevant to user's query
- Input: Product x
- Action: Score the relevance of x to user's query
- Feature Map:

$$\phi(x) = [1, N(x)],$$

where $N(x)$ = number of people who bought x .

- We expect a monotonic relationship between $N(x)$ and relevance, but...

Saturation: The Issue

Is relevance linear in $N(x)$?

- Relevance score reflects confidence in relevance prediction.
 - Are we 10 times more confident if $N(x) = 1000$ vs $N(x) = 100$?
-
- Bigger is better... but not that much better.

Saturation: Solve with nonlinear transform

- Smooth nonlinear transformation:

$$\phi(x) = [1, \log\{1 + N(x)\}]$$

- $\log(\cdot)$ good for values with large dynamic ranges
- *Does it matter what base we use in the log?*

From Percy Liang's "Lecture 3" slides from Stanford's CS221, Autumn 2014.

Saturation: Solve by discretization

- Discretization (a discontinuous transformation):

$$\phi(x) = (1(0 \leq N(x) < 10), 1(10 \leq N(x) < 100), \dots)$$

- Small buckets allow quite flexible relationship

Interactions: The Issue

- Input: Patient information x
- Action: Health score $y \in \mathbf{R}$ (higher is better)
- Feature Map

$$\phi(x) = [\text{height}(x), \text{weight}(x)]$$

- Issue: It's the weight **relative** to the height that's important.
- Impossible to get with these features and a linear classifier.
- Need some **interaction** between height and weight.

From Percy Liang's "Lecture 3" slides from Stanford's CS221, Autumn 2014.

Interactions: Approach 1

- Google “ideal weight from height”
- J. D. Robinson’s “ideal weight” formula (for a male):

$$\text{weight}(\text{kg}) = 52 + 1.9 [\text{height}(\text{in}) - 60]$$

- Make score square deviation between $\text{height}(h)$ and ideal weight(w)

$$f(x) = (52 + 1.9 [h(x) - 60] - w(x))^2$$

- WolframAlpha for complicated Mathematics:

$$f(x) = 3.61h(x)^2 - 3.8h(x)w(x) - 235.6h(x) + w(x)^2 + 124w(x) + 3844$$

Interactions: Approach 2

- Just include all second order features:

$$\phi(x) = \left[1, h(x), w(x), h(x)^2, w(x)^2, \underbrace{h(x)w(x)}_{\text{cross term}} \right]$$

- More flexible, no Google, no WolframAlpha.

General Principle

Simpler building blocks replace a single “smart” feature.

Predicate Features and Interaction Terms

Definition

A **predicate** on the input space \mathcal{X} is a function $P : \mathcal{X} \rightarrow \{\text{True}, \text{False}\}$.

- Many features take this form:
 - $x \mapsto s(x) = 1$ (subject is sleeping)
 - $x \mapsto d(x) = 1$ (subject is driving)
- For predicates, interaction terms correspond to **AND** conjunctions:
 - $x \mapsto s(x)d(x) = 1$ (subject is sleeping AND subject is driving)

Example: Monomial Interaction Terms

- Suppose we start with $x = (1, x_1, \dots, x_d) \in \mathbf{R}^{d+1} = \mathcal{X}$.
- To get a more expressive hypothesis space, we want to add interaction terms.
- Consider adding all monomials of degree M : $x_1^{p_1} \dots x_d^{p_d}$, with $p_1 + \dots + p_d = M$.
- How many features will we end up with?
- $\binom{M+d-1}{M}$ (“flower shop problem” from combinatorics)
- For $d = 40$ and $M = 8$, we get 314457495 features.
- That will make some extremely large data matrices...

Very large feature spaces have two potential issues:

- ① Overfitting
 - ② Memory and computational costs
- Overfitting we handle with regularization.
 - “**Kernel methods**” can (sometimes) help with memory and computational costs.

Kernel Methods: Motivation

SVM with Explicit Feature Map

- Let $\psi : \mathcal{X} \rightarrow \mathbf{R}^d$ be a feature map.
- The SVM optimization problem (with explicit feature map):

$$\min_{w \in \mathbf{R}^d} \frac{1}{2} \|w\|^2 + \frac{c}{n} \sum_{i=1}^n \max(0, 1 - y_i w^T \psi(x_i)).$$

- Last time we mentioned an equivalent optimization problem from Lagrangian duality...

SVM Dual Problem

- By Lagrangian duality, it is equivalent to solve the following optimization problem:

$$\begin{aligned} \max_{\alpha \in \mathbf{R}^n} \quad & \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i,j=1}^n \alpha_i \alpha_j y_i y_j \psi(x_j)^T \psi(x_i) \\ \text{s.t.} \quad & \sum_{i=1}^n \alpha_i y_i = 0 \quad \text{and} \quad \alpha_i \in \left[0, \frac{C}{n}\right] \quad i = 1, \dots, n. \end{aligned}$$

- If α^* is an optimal value, then

$$w^* = \sum_{i=1}^n \alpha_i^* y_i \psi(x_i) \quad \text{and} \quad \hat{f}(x) = \sum_{i=1}^n \alpha_i^* y_i \psi(x_i)^T \psi(x).$$

- Notice: $\psi(x)$ only shows up in an inner products with another $\psi(x')$.

Some Methods Can Be “Kernelized”

Definition

A method is **kernelized** if every feature vector $\psi(x)$ only appears inside an inner product with another feature vector $\psi(x')$. This applies to both the optimization problem and the prediction function.

- The SVM Dual is a kernelization of the original SVM formulation.
- We'll now introduce some special notation for these inner products $\langle \psi(x), \psi(x') \rangle \dots$

The Kernel Function

- **Input space:** \mathcal{X}
- **Feature space:** \mathcal{H} (a Hilbert space, i.e. an inner product space with projections, e.g. \mathbb{R}^d)
- **Feature map:** $\psi : \mathcal{X} \rightarrow \mathcal{H}$
- The **kernel function** corresponding to ψ is

$$k(x, x') = \langle \psi(x), \psi(x') \rangle,$$

where $\langle \cdot, \cdot \rangle$ is the inner product associated with \mathcal{H} .

The Kernel Function: Why do we need this?

- **Feature map:** $\psi : \mathcal{X} \rightarrow \mathcal{H}$
- The **kernel function** corresponding to ψ is

$$k(x, x') = \langle \psi(x), \psi(x') \rangle.$$

- Why introduce this new notation $k(x, x')$?
- We can often evaluate $k(x, x')$ without explicitly computing $\psi(x)$ and $\psi(x')$.
- For large feature spaces, can be much faster.

Kernel Evaluation Can Be Fast

Example

Quadratic feature map for $x = (x_1, \dots, x_d) \in \mathbf{R}^d$.

$$\psi(x) = (x_1, \dots, x_d, x_1^2, \dots, x_d^2, \sqrt{2}x_1x_2, \dots, \sqrt{2}x_i x_j, \dots, \sqrt{2}x_{d-1}x_d)^T$$

has dimension $O(d^2)$, but for any $x, x' \in \mathbf{R}^d$ and the standard Euclidean dot products,

$$k(x, x') = \langle \psi(x), \psi(x') \rangle = \langle x, x' \rangle + \langle x, x' \rangle^2$$

- Explicit computation of $k(x, x')$: $O(d^2)$
- Implicit computation of $k(x, x')$: $O(d)$

Kernels as Similarity Scores

- Often useful to think of the kernel function as a **similarity score**.
- But this is not a mathematically precise statement.
- There are many ways to design a similarity score.
- We will use kernel functions that correspond to inner products in some feature space.
- These are called **Mercer kernels**.

What are the Benefits of Kernelization?

- 1 Computational (when optimizing over \mathbf{R}^n is better than over \mathbf{R}^d)).
- 2 Can sometimes avoid any $O(d)$ operations
 - allows access to **infinite-dimensional feature spaces**.
- 3 Allows thinking in terms of “similarity” rather than features.

The Kernel Matrix

Definition

The **kernel matrix** for a kernel k on $x_1, \dots, x_n \in \mathcal{X}$ is

$$K = (k(x_i, x_j))_{i,j} = \begin{pmatrix} k(x_1, x_1) & \cdots & k(x_1, x_n) \\ \vdots & \ddots & \cdots \\ k(x_n, x_1) & \cdots & k(x_n, x_n) \end{pmatrix} \in \mathbf{R}^{n \times n}.$$

- In ML this is also called a **Gram matrix**, but traditionally (in linear algebra),
 - Gram matrices are defined without reference to a kernel or feature map.

The Kernel Matrix

- The kernel matrix summarizes all the information we need about the training inputs x_1, \dots, x_n to solve a kernelized optimization problem.
- e.g. in the kernelized SVM, we can replace $\psi(x_i)^T \psi(x_j)$ with K_{ij} :

$$\begin{aligned} \sup_{\alpha} \quad & \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i,j=1}^n \alpha_i \alpha_j y_i y_j K_{ij} \\ \text{s.t.} \quad & \sum_{i=1}^n \alpha_i y_i = 0 \quad \text{and} \quad \alpha_i \in \left[0, \frac{c}{n}\right] \quad i = 1, \dots, n. \end{aligned}$$

The “Kernel Trick”

- 1 Given a kernelized ML algorithm (i.e. all $\psi(x)$'s show up as $\langle \psi(x), \psi(x') \rangle$).
- 2 Can swap out the inner product for a new kernel function.
- 3 New kernel may correspond to a very high-dimensional feature space.
- 4 Once the kernel matrix is computed, the computational cost depends on number of data points, rather than the dimension of feature space.

The **trick** is that once you've implemented your method in terms of a kernel matrix, you can go from a kernel corresponding to a very small feature vector to a kernel corresponding to a very large (even infinite dimensional) feature vector, without changing your code, just by swapping one kernel matrix for another. Runtime is unaffected, after the kernel matrix is computed.

Our Plan

- Present our principal tool for kernelization: the **representer theorem**
- To keep things clean, we'll drop the explicit feature map until we need it: $\psi(x) = x$.
- Discuss specific cases of kernel ridge regression and kernel SVM
- Discuss several kernels, including the famous RBF kernel.
- Discuss how to create a kernel without an explicit feature map.

The Representer Theorem to Kernelize

The Representer Theorem

Theorem (Representer Theorem)

Let

$$J(w) = R(\|w\|) + L(\langle w, x_1 \rangle, \dots, \langle w, x_n \rangle),$$

where

- $w, x_1, \dots, x_n \in \mathcal{H}$ for some Hilbert space \mathcal{H} . (We typically have $\mathcal{H} = \mathbf{R}^d$.)
- $\|\cdot\|$ is the norm corresponding to the inner product of \mathcal{H} . (i.e. $\|w\| = \sqrt{\langle w, w \rangle}$)
- $R: [0, \infty) \rightarrow \mathbf{R}$ is nondecreasing (**Regularization term**), and
- $L: \mathbf{R}^n \rightarrow \mathbf{R}$ is arbitrary (**Loss term**).

If $J(w)$ has a minimizer, then it **has a minimizer of the form** $w^* = \sum_{i=1}^n \alpha_i x_i$.

[If R is strictly increasing, then all minimizers have this form. (Proof in homework.)]

Rewriting the Objective Function

- Define the training score function $s : \mathbf{R}^d \rightarrow \mathbf{R}^n$ by

$$s(w) = \begin{pmatrix} \langle w, x_1 \rangle \\ \vdots \\ \langle w, x_n \rangle \end{pmatrix},$$

which gives the **training score vector** for any w .

- We can then rewrite the objective function as

$$J(w) = R(\|w\|) + L(s(w)),$$

where now $L : \mathbf{R}^{n \times 1} \rightarrow \mathbf{R}$ takes a column vector as input.

- This will allow us to have a slick reparametrized version...

Reparametrize the Generalized Objective

- By the Representer Theorem, it's sufficient to minimize $J(w)$ for w of the form $\sum_{i=1}^n \alpha_i x_i$.
- Plugging this form into $J(w)$, we see we can just minimize

$$J_0(\alpha) = R\left(\left\|\sum_{i=1}^n \alpha_i x_i\right\|\right) + L\left(s\left(\sum_{i=1}^n \alpha_i x_i\right)\right)$$

over $\alpha = (\alpha_1, \dots, \alpha_n)^T \in \mathbf{R}^{n \times 1}$.

- With some new notation, we can substantially simplify
 - the norm piece $\|w\| = \left\|\sum_{i=1}^n \alpha_i x_i\right\|$, and
 - the score piece $s(w) = s\left(\sum_{i=1}^n \alpha_i x_i\right)$.

Simplifying the Reparametrized Norm

- For the norm piece $\|w\| = \|\sum_{i=1}^n \alpha_i x_i\|$, we have

$$\begin{aligned}\|w\|^2 &= \langle w, w \rangle \\ &= \left\langle \sum_{i=1}^n \alpha_i x_i, \sum_{j=1}^n \alpha_j x_j \right\rangle \\ &= \sum_{i,j=1}^n \alpha_i \alpha_j \langle x_i, x_j \rangle.\end{aligned}$$

- This expression involves the n^2 inner products between all pairs of input vectors.
- We often put those values together into a matrix...

The Gram Matrix

Definition

The **Gram matrix** of a set of points x_1, \dots, x_n in an inner product space is defined as

$$K = (\langle x_i, x_j \rangle)_{i,j} = \begin{pmatrix} \langle x_1, x_1 \rangle & \cdots & \langle x_1, x_n \rangle \\ \vdots & \ddots & \cdots \\ \langle x_n, x_1 \rangle & \cdots & \langle x_n, x_n \rangle \end{pmatrix}.$$

- This is the traditional definition from linear algebra.
- The Gram matrix is a special case of a **kernel matrix** for the identity feature map.
- That's why we write K for the Gram matrix instead of G , as done in elsewhere.
- NOTE: In ML, we often use Gram matrix and kernel matrix to mean the same thing. Don't get too hung up on the definitions.

Example: Gram Matrix for the Dot Product

- Consider $x_1, \dots, x_n \in \mathbf{R}^{d \times 1}$ with the standard inner product $\langle x, x' \rangle = x^T x'$.
- Let $X \in \mathbf{R}^{n \times d}$ be the **design matrix**, which has each input vector as a row:

$$X = \begin{pmatrix} -x_1^T - \\ \vdots \\ -x_n^T - \end{pmatrix}.$$

- Then the Gram matrix is

$$\begin{aligned} K &= \begin{pmatrix} x_1^T x_1 & \cdots & x_1^T x_n \\ \vdots & \ddots & \vdots \\ x_n^T x_1 & \cdots & x_n^T x_n \end{pmatrix} = \begin{pmatrix} -x_1^T - \\ \vdots \\ -x_n^T - \end{pmatrix} \begin{pmatrix} | & \cdots & | \\ x_1 & \cdots & x_n \\ | & \cdots & | \end{pmatrix} \\ &= XX^T \end{aligned}$$

Simplifying the Reparametrized Norm

- With $w = \sum_{i=1}^n \alpha_i x_i$, we have

$$\begin{aligned}\|w\|^2 &= \langle w, w \rangle \\ &= \left\langle \sum_{i=1}^n \alpha_i x_i, \sum_{j=1}^n \alpha_j x_j \right\rangle \\ &= \sum_{i,j=1}^n \alpha_i \alpha_j \langle x_i, x_j \rangle \\ &= \alpha^T K \alpha.\end{aligned}$$

Simplifying the Training Score Vector

- The score for x_j for $w = \sum_{i=1}^n \alpha_i x_i$ is

$$\langle w, x_j \rangle = \left\langle \sum_{i=1}^n \alpha_i x_i, x_j \right\rangle = \sum_{i=1}^n \alpha_i \langle x_i, x_j \rangle$$

- The training score vector is

$$\begin{aligned} s \left(\sum_{i=1}^n \alpha_i x_i \right) &= \begin{pmatrix} \sum_{i=1}^n \alpha_i \langle x_i, x_1 \rangle \\ \vdots \\ \sum_{i=1}^n \alpha_i \langle x_i, x_n \rangle \end{pmatrix} = \begin{pmatrix} \alpha_1 \langle x_1, x_1 \rangle + \cdots + \alpha_n \langle x_n, x_1 \rangle \\ \vdots \\ \alpha_1 \langle x_1, x_n \rangle + \cdots + \alpha_n \langle x_n, x_n \rangle \end{pmatrix} \\ &= \begin{pmatrix} \langle x_1, x_1 \rangle & \cdots & \langle x_1, x_n \rangle \\ \vdots & \ddots & \cdots \\ \langle x_n, x_1 \rangle & \cdots & \langle x_n, x_n \rangle \end{pmatrix} \begin{pmatrix} \alpha_1 \\ \vdots \\ \alpha_n \end{pmatrix} \\ &= K \alpha \end{aligned}$$

Reparametrized Objective

- Putting it all together, our reparametrized objective function can be written as

$$\begin{aligned} J_0(\alpha) &= R\left(\left\|\sum_{i=1}^n \alpha_i x_i\right\|\right) + L\left(s\left(\sum_{i=1}^n \alpha_i x_i\right)\right) \\ &= R\left(\sqrt{\alpha^T K \alpha}\right) + L(K\alpha), \end{aligned}$$

which we minimize over $\alpha \in \mathbf{R}^n$.

- All information needed about x_1, \dots, x_n is summarized in the Gram matrix K .
- We're now minimizing over \mathbf{R}^n rather than \mathbf{R}^d .
- If $d \gg n$, this can be a big win computationally (at least once K is computed).

Reparametrizing Predictions

- Suppose we've found

$$\alpha^* \in \arg \min_{\alpha \in \mathbf{R}^n} R\left(\sqrt{\alpha^T K \alpha}\right) + L(K\alpha).$$

- Then we know $w^* = \sum_{i=1}^n \alpha_i^* x_i$ satisfies

$$w^* \in \arg \min_{w \in \mathcal{H}} R(\|w\|) + L(\langle w, x_1 \rangle, \dots, \langle w, x_n \rangle).$$

- The prediction on a new point $x \in \mathcal{H}$ is

$$\hat{f}(x) = \langle w^*, x \rangle = \sum_{i=1}^n \alpha_i^* \langle x_i, x \rangle.$$

- To make a new prediction, we may need to touch all the training inputs x_1, \dots, x_n .

- It will be convenient to define the following column vector for any $x \in \mathcal{H}$:

$$k_x = \begin{pmatrix} \langle x_1, x \rangle \\ \vdots \\ \langle x_n, x \rangle \end{pmatrix}$$

- Then we can write our predictions on a new point x as

$$\hat{f}(x) = k_x^T \alpha^*$$

Summary So Far

- Original plan:
 - Find $w^* \in \arg \min_{w \in \mathcal{H}} R(\|w\|) + L(\langle w, x_1 \rangle, \dots, \langle w, x_n \rangle)$
 - Predict with $\hat{f}(x) = \langle w^*, x \rangle$.
- We showed that the following is equivalent:
 - Find $\alpha^* \in \arg \min_{\alpha \in \mathbb{R}^n} R(\sqrt{\alpha^T K \alpha}) + L(K \alpha)$
 - Predict with $\hat{f}(x) = k_x^T \alpha^*$, where

$$K = \begin{pmatrix} \langle x_1, x_1 \rangle & \cdots & \langle x_1, x_n \rangle \\ \vdots & \ddots & \dots \\ \langle x_n, x_1 \rangle & \cdots & \langle x_n, x_n \rangle \end{pmatrix} \quad \text{and} \quad k_x = \begin{pmatrix} \langle x_1, x \rangle \\ \vdots \\ \langle x_n, x \rangle \end{pmatrix}$$

- Every element $x \in \mathcal{H}$ occurs inside an inner products with a training input $x_i \in \mathcal{H}$.

Definition

A method is **kernelized** if every feature vector $\psi(x)$ only appears inside an inner product with another feature vector $\psi(x')$. This applies to both the optimization problem and the prediction function.

- Here we are using $\psi(x) = x$. Thus finding

$$\alpha^* \in \arg \min_{\alpha \in \mathbf{R}^n} R\left(\sqrt{\alpha^T K \alpha}\right) + L(K\alpha)$$

and making predictions with $\hat{f}(x) = k_x^T \alpha^*$ is a **kernelization** of finding

$$w^* \in \arg \min_{w \in \mathcal{H}} R(\|w\|) + L(\langle w, x_1 \rangle, \dots, \langle w, x_n \rangle)$$

and making predictions with $\hat{f}(x) = \langle w^*, x \rangle$.

- Once we have kernelized:
 - $\alpha^* \in \arg \min_{\alpha \in \mathbb{R}^n} R\left(\sqrt{\alpha^T K \alpha}\right) + L(K\alpha)$
 - $\hat{f}(x) = k_x^T \alpha^*$
- We can do the “kernel trick”.
- Replace each $\langle x, x' \rangle$ by $k(x, x')$, for any kernel function k , where $k(x, x') = \langle \psi(x), \psi(x') \rangle$.
- Predictions

$$\hat{f}(x) = \sum_{i=1}^n \alpha_i^* k(x_i, x)$$

Kernel Ridge Regression

Kernelizing Ridge Regression

- Ridge Regression:

$$\min_{w \in \mathbf{R}^d} \frac{1}{n} \|Xw - y\|^2 + \lambda \|w\|^2$$

- Plugging in $w = \sum_{i=1}^n \alpha_i x_i$, we get the kernelized ridge regression objective function:

$$\min_{\alpha \in \mathbf{R}^n} \frac{1}{n} \|K\alpha - y\|^2 + \lambda \alpha^T K \alpha$$

- This is usually just called **kernel ridge regression**.

Kernel Ridge Regression Solutions

- For $\lambda > 0$, the **ridge regression solution** is

$$w^* = (X^T X + \lambda I)^{-1} X^T y$$

- and the **kernel ridge regression solution** is

$$\begin{aligned} \alpha^* &= (X X^T + \lambda I)^{-1} y \\ &= (K + \lambda I)^{-1} y \end{aligned}$$

- (Shown in homework.)
- For ridge regression we're dealing with a $d \times d$ matrix.
- For kernel ridge regression we're dealing an $n \times n$ matrix.

Predictions

- Predictions in terms of w^* :

$$\hat{f}(x) = x^T w^*$$

- Predictions in terms of α^* :

$$\hat{f}(x) = k_x^T \alpha^* = \sum_{i=1}^n \alpha_i^* x_i^T x$$

- For kernel ridge regression, need to access all training inputs x_1, \dots, x_n to predict.
- For SVM, we may not...

Kernel SVM

Kernelized SVM (From Representer Theorem)

- The SVM objective:

$$\min_{w \in \mathbf{R}^d} \frac{1}{2} \|w\|^2 + \frac{c}{n} \sum_{i=1}^n \max(0, 1 - y_i w^T x_i).$$

- Plugging in $w = \sum_{i=1}^n \alpha_i x_i$, we get

$$\min_{\alpha \in \mathbf{R}^n} \frac{1}{2} \alpha^T K \alpha + \frac{c}{n} \sum_{i=1}^n \max(0, 1 - y_i (K \alpha)_i)$$

- Predictions with

$$\hat{f}(x) = x^T w^* = \sum_{i=1}^n \alpha_i^* x_i^T x.$$

- This is one way to kernelize SVM...

Kernelized SVM (From Lagrangian Duality)

- Kernelized SVM from computing the Lagrangian Dual Problem:

$$\begin{aligned} \max_{\alpha \in \mathbf{R}^n} \quad & \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i,j=1}^n \alpha_i \alpha_j y_i y_j x_j^T x_i \\ \text{s.t.} \quad & \sum_{i=1}^n \alpha_i y_i = 0 \\ & \alpha_i \in \left[0, \frac{C}{n}\right] \quad i = 1, \dots, n. \end{aligned}$$

- If α^* is an optimal value, then

$$w^* = \sum_{i=1}^n \alpha_i^* y_i x_i \quad \text{and} \quad \hat{f}(x) = \sum_{i=1}^n \alpha_i^* y_i x_i^T x.$$

- Note that the prediction function is also kernelized.

Sparsity in the Data from Complementary Slackness

- Kernelized predictions given by

$$\hat{f}(x) = \sum_{i=1}^n \alpha_i^* y_i x_i^T x.$$

- By a Lagrangian duality analysis (specifically from complementary slackness), we find

$$y_i \hat{f}(x_i) < 1 \implies \alpha_i^* = \frac{c}{n}$$

$$y_i \hat{f}(x_i) = 1 \implies \alpha_i^* \in \left[0, \frac{c}{n}\right]$$

$$y_i \hat{f}(x_i) > 1 \implies \alpha_i^* = 0$$

- So we can leave out any x_i “on the good side of the margin” ($y_i \hat{f}(x_i) > 1$).
- x_i 's that we must keep, because $\alpha_i^* \neq 0$, are called **support vectors**.

Kernels

Linear Kernel

- Input space: $\mathcal{X} = \mathbf{R}^d$
- Feature space: $\mathcal{H} = \mathbf{R}^d$, with standard inner product
- Feature map

$$\psi(x) = x$$

- Kernel:

$$k(x, x') = x^T x'$$

Quadratic Kernel in \mathbf{R}^d

- Input space $\mathcal{X} = \mathbf{R}^d$
- Feature space: $\mathcal{H} = \mathbf{R}^D$, where $D = d + \binom{d}{2} \approx d^2/2$.
- Feature map:

$$\psi(x) = (x_1, \dots, x_d, x_1^2, \dots, x_d^2, \sqrt{2}x_1x_2, \dots, \sqrt{2}x_ix_j, \dots, \sqrt{2}x_{d-1}x_d)^T$$

- Then for $\forall x, x' \in \mathbf{R}^d$

$$\begin{aligned}k(x, x') &= \langle \psi(x), \psi(x') \rangle \\ &= \langle x, x' \rangle + \langle x, x' \rangle^2\end{aligned}$$

- Computation for inner product with explicit mapping: $O(d^2)$
- Computation for implicit kernel calculation: $O(d)$.

Polynomial Kernel in \mathbf{R}^d

- Input space $\mathcal{X} = \mathbf{R}^d$
- Kernel function:

$$k(x, x') = (1 + \langle x, x' \rangle)^M$$

- Corresponds to a feature map with all monomials up to degree M .
- For any M , computing the kernel has same computational cost
- Cost of explicit inner product computation grows rapidly in M .

The RBF Kernel

Radial Basis Function (RBF) / Gaussian Kernel

- Input space $\mathcal{X} = \mathbf{R}^d$

$$k(x, x') = \exp\left(-\frac{\|x - x'\|^2}{2\sigma^2}\right),$$

where σ^2 is known as the bandwidth parameter.

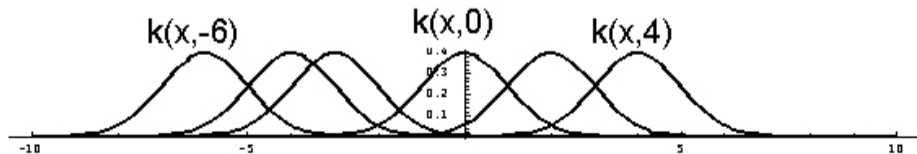
- Does it act like a similarity score?
- Why “radial”?
- Have we departed from our “inner product of feature vector” recipe?
 - Yes and no: corresponds to an infinite dimensional feature vector
- Probably the most common nonlinear kernel.

RBF Basis

- Input space $\mathcal{X} = \mathbb{R}$
- Output space: $\mathcal{Y} = \mathbb{R}$
- RBF kernel $k(w, x) = \exp\left(- (w - x)^2\right)$.
- Suppose we have 6 training examples: $x_i \in \{-6, -4, -3, 0, 2, 4\}$.
- If representer theorem applies, then

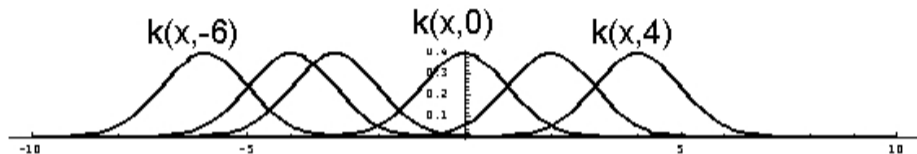
$$f(x) = \sum_{i=1}^6 \alpha_i k(x_i, x).$$

- f is a linear combination of 6 basis functions of form $k(x_i, \cdot)$:

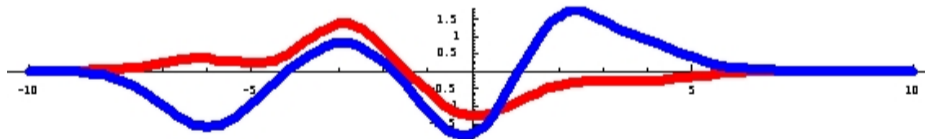


RBF Predictions

- Basis functions



- Predictions of the form $f(x) = \sum_{i=1}^6 \alpha_i k(x_i, x)$:



- When kernelizing with RBF kernel, prediction functions always look this way.
- (Whether we get w from SVM, ridge regression, etc...)

RBF Feature Space: The Sequence Space ℓ_2

- To work with infinite dimensional feature vectors, we need a space with certain properties.
 - an inner product
 - a norm related to the inner product
 - projection theorem: $x = x_{\perp} + x_{\parallel}$ where $x_{\parallel} \in S = \text{span}(w_1, \dots, w_n)$ and $\langle x_{\perp}, s \rangle = 0 \quad \forall s \in S$.
- Basically, we need a Hilbert space.

Definition

ℓ_2 is the space of all real-valued sequences: $(x_0, x_1, x_2, x_3, \dots)$ with $\sum_{i=0}^{\infty} x_i^2 < \infty$.

Theorem

*With the inner product $\langle x, x' \rangle = \sum_{i=0}^{\infty} x_i x'_i$, ℓ_2 is a **Hilbert space**.*

The Infinite Dimensional Feature Vector for RBF

- Consider RBF kernel (1-dim): $k(x, x') = \exp\left(-\frac{(x - x')^2}{2}\right)$
- We claim that $\psi : \mathbf{R} \rightarrow \ell_2$, defined by

$$[\psi(x)]_j = \frac{1}{\sqrt{j!}} e^{-x^2/2} x^j$$

gives the “infinite-dimensional feature vector” corresponding to RBF kernel.

- Is this mapping even well-defined? Is $\psi(x)$ even an element of ℓ_2 ?
- Yes:

$$\sum_{j=0}^{\infty} \frac{1}{j!} e^{-x^2} x^{2j} = e^{-x^2} \sum_{j=0}^{\infty} \frac{(x^2)^j}{j!} = 1 < \infty$$

The Infinite Dimensional Feature Vector for RBF

- Does feature vector $[\psi(x)]_n = \frac{1}{\sqrt{j!}} e^{-x^2/2} x^j$ actually correspond to the RBF kernel?
- Yes! Proof:

$$\begin{aligned}\langle \psi(x), \psi(x') \rangle &= \sum_{j=0}^{\infty} \frac{1}{j!} e^{-(x^2+(x')^2)/2} x^j (x')^j \\ &= e^{-(x^2+(x')^2)/2} \sum_{j=0}^{\infty} \frac{(xx')^j}{j!} \\ &= \exp\left(-\left[x^2 + (x')^2\right]/2\right) \exp(xx') \\ &= \exp\left(-\left[(x-x')^2/2\right]\right)\end{aligned}$$

QED

When is $k(x, x')$ a kernel function? (Mercer's Theorem)

How to Get Kernels?

- 1 Explicitly construct $\psi(x) : \mathcal{X} \rightarrow \mathbf{R}^d$ and define $k(x, x') = \psi(x)^T \psi(x')$.
- 2 Directly define the kernel function $k(x, x')$, and verify it corresponds to $\langle \psi(x), \psi(x') \rangle$ for some ψ .

There are many theorems to help us with the second approach

Positive Semidefinite Matrices

Definition

A real, symmetric matrix $M \in \mathbf{R}^{n \times n}$ is **positive semidefinite (psd)** if for any $x \in \mathbf{R}^n$,

$$x^T M x \geq 0.$$

Theorem

The following conditions are each necessary and sufficient for a symmetric matrix M to be positive semidefinite:

- *M can be factorized as $M = R^T R$, for some matrix R .*
- *All eigenvalues of M are greater than or equal to 0.*

Positive Semidefinite Function

Definition

A symmetric kernel function $k : \mathcal{X} \times \mathcal{X} \rightarrow \mathbf{R}$ is **positive semidefinite (psd)** if for any finite set $\{x_1, \dots, x_n\} \in \mathcal{X}$, the kernel matrix on this set

$$K = (k(x_i, x_j))_{i,j} = \begin{pmatrix} k(x_1, x_1) & \cdots & k(x_1, x_n) \\ \vdots & \ddots & \cdots \\ k(x_n, x_1) & \cdots & k(x_n, x_n) \end{pmatrix}$$

is a positive semidefinite matrix.

Mercer's Theorem

Theorem

A symmetric function $k(x, x')$ can be expressed as an inner product

$$k(x, x') = \langle \psi(x), \psi(x') \rangle$$

for some ψ if and only if $k(x, x')$ is **positive semidefinite**.

Generating New Kernels from Old

- Suppose $k, k_1, k_2 : \mathcal{X} \times \mathcal{X} \rightarrow \mathbf{R}$ are psd kernels. Then so are the following:

$$k_{\text{new}}(x, x') = k_1(x, x') + k_2(x, x')$$

$$k_{\text{new}}(x, x') = \alpha k(x, x')$$

$$k_{\text{new}}(x, x') = f(x)f(x') \text{ for any function } f(\cdot)$$

$$k_{\text{new}}(x, x') = k_1(x, x')k_2(x, x')$$

- See Appendix for details.
- Lots more theorems to help you construct new kernels from old...

Details on New Kernels from Old [Optional]

- Suppose k_1 and k_2 are psd kernels with feature maps ϕ_1 and ϕ_2 , respectively.
- Then

$$k_1(x, x') + k_2(x, x')$$

is a psd kernel.

- Proof: Concatenate the feature vectors to get

$$\phi(x) = (\phi_1(x), \phi_2(x)).$$

Then ϕ is a feature map for $k_1 + k_2$.

Closure under Positive Scaling

- Suppose k is a psd kernel with feature maps ϕ .
- Then for any $\alpha > 0$,

$$\alpha k$$

is a psd kernel.

- Proof: Note that

$$\phi(x) = \sqrt{\alpha}\phi(x)$$

is a feature map for αk .

Scalar Function Gives a Kernel

- For any function $f(x)$,

$$k(x, x') = f(x)f(x')$$

is a kernel.

- Proof: Let $f(x)$ be the feature mapping. (It maps into a 1-dimensional feature space.)

$$\langle f(x), f(x') \rangle = f(x)f(x') = k(x, x').$$

Closure under Hadamard Products

- Suppose k_1 and k_2 are psd kernels with feature maps ϕ_1 and ϕ_2 , respectively.
- Then

$$k_1(x, x') k_2(x, x')$$

is a psd kernel.

- Proof: Take the outer product of the feature vectors:

$$\phi(x) = \phi_1(x) [\phi_2(x)]^T.$$

Note that $\phi(x)$ is a matrix.

- Continued...

- Then

$$\begin{aligned}\langle \phi(x), \phi(x') \rangle &= \sum_{ij} \phi(x) \phi(x') \\ &= \sum_{ij} [\phi_1(x) [\phi_2(x)]^T]_{ij} [\phi_1(x') [\phi_2(x')]^T]_{ij} \\ &= \sum_{ij} [\phi_1(x)]_i [\phi_2(x)]_j [\phi_1(x')]_i [\phi_2(x')]_j \\ &= \left(\sum_i [\phi_1(x)]_i [\phi_1(x')]_i \right) \left(\sum_j [\phi_2(x)]_j [\phi_2(x')]_j \right) \\ &= k_1(x, x') k_2(x, x')\end{aligned}$$