## Tutorial 5 – Kernels and Gaussian Processes

CSC2541 Neural Net Training Dynamics - Winter 2022

Slides adapted from CSC2541: Scalable and Flexible Models of Uncertainty - Fall 2017

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## Recap: Basis Functions

- Basis functions allow us to use non-linear feature transformations.
- We can specify them by hand (examples below), or learn them automatically using a neural network.

$$\phi_{j}(x) = x^{j}$$

$$0.5$$

$$0$$

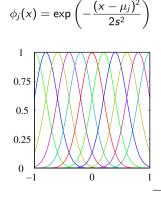
$$-0.5$$

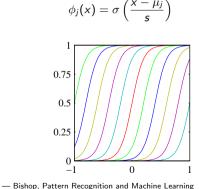
$$-1$$

$$-1$$

$$0$$

$$1$$





## Recap: Basis Functions

• How is this useful? We can use linear methods on non-linear features to yield non-linear decision boundaries and regression curves.

$$\phi\left(\begin{bmatrix}x_1\\x_2\end{bmatrix}\right) = \begin{bmatrix}x_1^2\\x_2^2\\\sqrt{2}x_1x_2\end{bmatrix}$$

— https://gregorygundersen.com/blog/2019/12/10/kernel-trick/

### Kernels: Motivation

### Generalized Linear Models (GLM)

- Fixed non-linear basis functions.
- Limited hypothesis space.
- Easy to optimize (convex).

### Neural Network (NN)

- Adaptive non-linear basis functions.
- Rich hypothesis space.
- Hard to optimize (non-convex).

#### Towards Kernel Methods

- Feature space in GLM and NN needs to be explicitly constructed.
- Can we use a large (possibly infinite) set of fixed non-linear basis functions without explicitly constructing this space?
- Yes, by using kernel methods!

#### Kernel Methods

- Kernel methods are instance-based learners: they assign a weight  $\theta_i$  to any training point  $\mathbf{x}_i$ .
- Predictions on new data points  $\mathbf{x}'$  make use of a kernel function  $\kappa(\cdot, \cdot)$  measuring the similarity of  $\mathbf{x}'$  with all points  $\mathbf{x}_i$  from the training set.
- Kernelized binary classification example:

$$\hat{y} = \operatorname{sgn} \sum_{i=1}^{n} \theta_{i} y_{i} \kappa(\mathbf{x}_{i}, \mathbf{x}')$$

where

- $y \in \{-1, +1\}$  is the label assigned to a data point **x**.
- $\theta_i$  is the weight for training example  $\mathbf{x}_i$ .
- $\kappa: \mathcal{X} \times \mathcal{X} \to \mathbb{R}$  is the kernel function measuring similarity between  $\mathbf{x}, \mathbf{x}' \in \mathbb{R}$ .

#### The Kernel Trick

- Let  $\phi(\cdot)$  be a set of not further specified basis functions mappings.
- Explicitly constructing a high-dimensional feature space is expensive.
- By using the kernel trick, we can implicitly perform operations in a high-dimensional feature space.
- In many algorithms, this feature space only appears as a dot product  $\langle \phi(\mathbf{x}), \phi(\mathbf{x}') \rangle = \phi(\mathbf{x})^{\top} \phi(\mathbf{x}')$  of input pairs  $\mathbf{x}, \mathbf{x}'$ .
- We define these dot products as the kernel function

$$\kappa(\mathbf{x}, \mathbf{x}') = \langle \phi(\mathbf{x}), \phi(\mathbf{x}') \rangle = \phi(\mathbf{x})^{\top} \phi(\mathbf{x}')$$

which can also be thought of as a similarity function between x and x'.

### **Dual Representation**

• Recall the regularized linear regression objective:

$$\mathcal{L}(\boldsymbol{\theta}) = \frac{1}{2} \sum_{n=1}^{N} (\boldsymbol{\theta}^{\top} \phi(\mathbf{x}_n) - y_n)^2 + \frac{\lambda}{2} \boldsymbol{\theta}^{\top} \boldsymbol{\theta}$$

• Finding optimal  $\theta$ :

$$\nabla_{\theta} \mathcal{L}(\theta) = \sum_{n=1}^{N} (\theta^{\top} \phi(\mathbf{x}_n) - y_n) \phi(\mathbf{x}_n) + \lambda \theta = 0$$
$$\theta = -\frac{1}{\lambda} \sum_{n=1}^{N} \underbrace{(\theta^{\top} \phi(\mathbf{x}_n) - y_n)}_{\bullet} \phi(\mathbf{x}_n)$$

• The weights  $\theta$  can be written as a linear combination of the training examples:

$$heta = \sum_{n=1}^N \mathbf{a}_n \phi(\mathbf{x}_n)$$
 where  $\mathbf{a} = \left[\mathbf{a}_1, \dots, \mathbf{a}_n \right]$  are called the dual parameters

## **Dual Representation**

• Substituting  $\theta$  back into linear regression  $y(\mathbf{x}) = \theta^{\top} \phi(\mathbf{x})$  yields:

$$\theta = \sum_{n=1}^{N} \mathbf{a}_n \phi(\mathbf{x}_n) \qquad y(\mathbf{x}) = \sum_{n=1}^{N} \mathbf{a}_n \phi(\mathbf{x}_n)^{\top} \phi(\mathbf{x}) = \sum_{n=1}^{N} \mathbf{a}_n \kappa(\mathbf{x}_n, \mathbf{x})$$

- The feature space only appears as a dot product.
- The kernel matrix, or gram matrix,  $\mathbf{K} \in \mathbb{R}^{N \times N}$  collects kernel values in a symmetric positive semi-definite matrix for all data points (Mercer's theorem):

$$\mathbf{K}_{ij} = \kappa(\mathbf{x}_i, \mathbf{x}_j) = \phi(\mathbf{x}_i)^{\top} \phi(\mathbf{x}_j)$$

• If a kernel defines such a kernel matrix, then the kernel is valid.

## Popular Kernels

### Polynomial Kernel

$$\kappa_{\text{Pol}}(\mathbf{x}, \mathbf{x}') = (\mathbf{x}^{\top} \mathbf{x}' + c)^{d}$$

$$0.5$$

$$0$$

$$-0.5$$

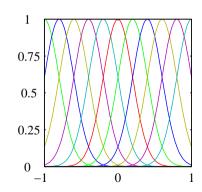
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$$0$$

# Squared Exponential Kernel

$$\kappa_{
m SE}(\mathbf{x},\mathbf{x}') = \sigma^2 \exp\left(-rac{(\mathbf{x}-\mathbf{x}')^2}{2\ell^2}
ight)$$



## Kernel Composition Rules

Let  $\kappa_1(\mathbf{x}, \mathbf{x}')$  and  $\kappa_2(\mathbf{x}, \mathbf{x}')$  be valid kernels, then the following kernels are also valid:

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = c\kappa_1(\mathbf{x}, \mathbf{x}') \quad \forall c > 0$$

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = f(\mathbf{x})\kappa_1(\mathbf{x}, \mathbf{x}')f(\mathbf{x}') \quad \forall$$

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = g(\kappa_1(\mathbf{x}, \mathbf{x}'))$$
  $g$  is polynomial with coefficients  $\geq 0$ .

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = \exp(\kappa_1(\mathbf{x}, \mathbf{x}'))$$

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = \kappa_1(\mathbf{x}, \mathbf{x}') + \kappa_2(\mathbf{x}, \mathbf{x}')$$
 kernel OR-ing

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = \kappa_1(\mathbf{x}, \mathbf{x}')\kappa_2(\mathbf{x}, \mathbf{x}')$$
 kernel AND-ing

• 
$$\kappa(\mathbf{x}, \mathbf{x}') = \mathbf{x}^{\top} \mathbf{A} \mathbf{x}'$$
 **A** symmetric and p.s.d.

Check out the Kernel Cookbook:

https://www.cs.toronto.edu/~duvenaud/cookbook/

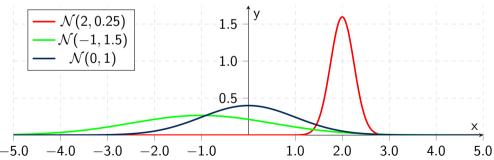


# Recap: Multivariate Gaussian

- Handy tool for Bayesian inference on real-valued variables
- General multivariate PDF:

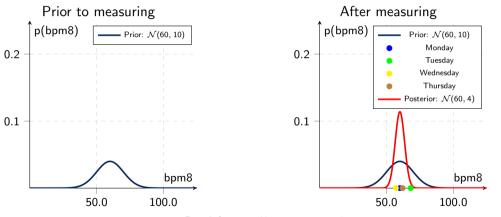
$$\mathbf{x} \sim \mathcal{N}_D(oldsymbol{\mu}, oldsymbol{\Sigma}) = rac{1}{\sqrt{(2\pi)^D |oldsymbol{\Sigma}|}} e^{-rac{1}{2}(\mathbf{x} - oldsymbol{\mu})^ op oldsymbol{\Sigma}^{-1}(\mathbf{x} - oldsymbol{\mu})}$$

ullet Some examples of D=1 Gaussians



# Bayesian Parameter Estimation Example

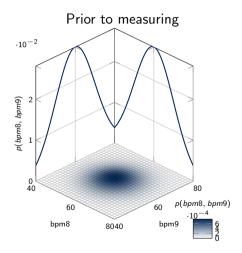
#### Measure your heart rate at 8am

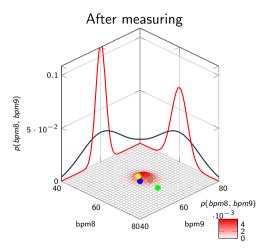


 $-- {\tt Example from http://videolectures.net/mlss2012\_cunningham\_gaussian\_processes/}$ 

# Bayesian Parameter Estimation Example

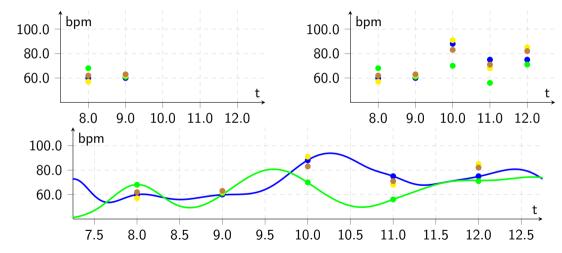
#### Measure your heart rate at 8am and 9am





# Bayesian Parameter Estimation Example

Measuring your heart rate throughout the day



## $\mathcal{GP}$ Definition

A Gaussian process describes a distribution over functions (infinitely long vectors).

- Notation:  $f(\mathbf{x}) \sim \mathcal{GP}(m(\mathbf{x}), \kappa(\mathbf{x}, \mathbf{x}))$
- Mean function:  $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$
- Covariance function:  $\kappa(\mathbf{x}, \mathbf{x}') = \mathbb{E}[(f(\mathbf{x}) m(\mathbf{x}))(f(\mathbf{x}') m(\mathbf{x}'))]$

We have data points  $\boldsymbol{X} = [\boldsymbol{x}_1^\top, \dots, \boldsymbol{x}_n^\top]^\top$  and are interested in their function values  $f(\boldsymbol{X}) = (f(\boldsymbol{x}_1), \dots, f(\boldsymbol{x}_n))^\top$ .

A Gaussian process is a collection of random variables, any finite number of which have joint Gaussian distribution.

f(x) is one such subset and has (prior) joint Gaussian distribution.

### GP Mean and Covariance

#### The mean function m

- The mean function  $m(\cdot)$  encodes the a-priori expectation of the function.
- m(x) will dominate the inference result in case we have not yet observed data similar to x.
- Typical choice: zero-centering the data:  $m(\mathbf{x}) = 0$

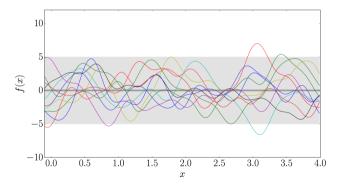
#### The covariance function $\kappa$

- $\kappa(\mathbf{x}, \mathbf{x}')$  measures similarity between  $\mathbf{x}$  and  $\mathbf{x}' \to \text{similar}$  data points have similar function values.
- κ is a Mercer kernel.
- Typical choice: squared exponential kernel:  $\kappa(\mathbf{x},\mathbf{x}') = \sigma^2 e^{-\frac{(\mathbf{x}-\mathbf{x}')^\top(\mathbf{x}-\mathbf{x}')}{2\ell^2}}$  where  $\sigma$ defines the height and  $\ell$  the width of the kernel.

## Drawing Samples From The Prior

Same procedure as for multivariate Gaussians:

- 1. Generate  $u \in \mathbb{R}^D$  by drawing d samples from  $\mathcal{N}(\mathbf{0}, \mathbf{I}_D)$ .
- 2. Perform Cholesky decomposition  $\Sigma = LL^{\top}$ .
- 3. Compute  $\mathbf{y} = \boldsymbol{\mu} + \boldsymbol{L}\boldsymbol{u}$  where  $\mathbf{y} \sim \mathcal{N}(\boldsymbol{\mu}, \boldsymbol{\Sigma})$ .



### The Joint Distribution

We have training data  $\mathbf{X} \in \mathbb{R}^{N \times D}$ , corresponding observations  $\mathbf{y} = f(\mathbf{X})$ , and test data points  $\mathbf{X}_* \in \mathbb{R}^{N_* \times D}$  for which we want to infer function values  $\mathbf{v}_* = f(\mathbf{X}_*)$ . The GP defines the following joint distribution

$$\rho(\mathbf{y}, \mathbf{y}_* | \mathbf{X}, \mathbf{X}_*) = \begin{pmatrix} \mathbf{y} \\ \mathbf{y}_* \end{pmatrix} \sim \mathcal{N} \left( \begin{bmatrix} m(\mathbf{X}) \\ m(\mathbf{X}_*) \end{bmatrix}, \begin{bmatrix} \mathbf{K} + \sigma_n^2 \mathbf{I} & \mathbf{K}_* \\ \mathbf{K}_*^\top & \mathbf{K}_{**} \end{bmatrix} \right)$$

where

$$\mathbf{K} = \kappa(\mathbf{X}, \mathbf{X})$$
  $\mathbf{K}_* = \kappa(\mathbf{X}, \mathbf{X}_*)$   $\mathbf{K}_{**} = \kappa(\mathbf{X}_*, \mathbf{X}_*).$ 

Typically, data points are corrupted by noise  $\rightarrow$  our functions should not act as interpolators. We therefore assume

$$y_i = f(\mathbf{x}_i) + \epsilon$$
 where  $\epsilon \sim \mathcal{N}(0, \sigma_n^2)$ .

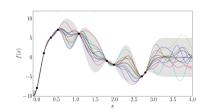
### Inference with Gaussian Processes

Inferring an unknown function value and its covariance follows from conditioning multivariate Gaussians:

$$ho( extbf{ extit{y}}_*| extbf{ extit{y}}, extbf{ extit{X}}, extbf{ extit{X}}_*) \sim \mathcal{N}(m{\mu}, m{\Sigma})$$

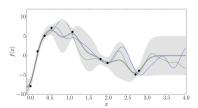
#### Non-noisy case

- $\mu = m(X_*) + K_*^{\top} K^{-1} (y m(X))$
- $\Sigma = K_{**} K_{*}^{\top} K^{-1} K_{*}$

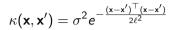


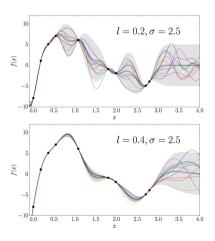
#### Noisy case

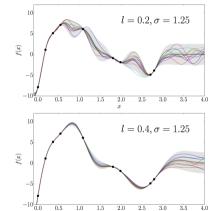
- $\mu = m(X_*) + K_*^{\top} (K + \sigma_n^2 I)^{-1} (y m(X))$
- $\bullet \ \ \boldsymbol{\Sigma} = \boldsymbol{\mathcal{K}}_{**} \boldsymbol{\mathcal{K}}_*^\top (\boldsymbol{\mathcal{K}} + \sigma_n^2 \boldsymbol{I})^{-1} \boldsymbol{\mathcal{K}}_*$



# Influence of Kernel Hyperparameters







### References I

#### Useful links

- https://distill.pub/2019/visual-exploration-gaussian-processes/
- http://www.infinitecuriosity.org/vizgp/
- https://mlg.eng.cam.ac.uk/tutorials/06/es.pdf
- https://xavierbourretsicotte.github.io/Kernel\_feature\_map.html
- https://www.youtube.com/watch?v=nzSBvINmg28
- https://www.youtube.com/watch?v=exqpaqaPG2M
- Christopher M Bishop and Nasser M Nasrabadi, *Pattern recognition and machine learning*, vol. 4, Springer, 2006.
- Christopher K Williams and Carl Edward Rasmussen, *Gaussian processes for machine learning*, vol. 2, MIT press Cambridge, MA, 2006.