# **Introduction to Machine Learning**

# **Gaussian Processes Covariance functions for GPs**

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#### **Learning goals**

- Covariance functions encode key assumptions about the GP
- Know common covariance functions like squared exponential and Matérn

# **COVARIANCE FUNCTION OF A GP**

The marginalization property of the Gaussian process implies that for any finite set of input values, the corresponding vector of function values is Gaussian:

$$
\textbf{\textit{f}}=\left[f\left(\textbf{x}^{(1)}\right),...,f\left(\textbf{x}^{(n)}\right)\right]\sim\mathcal{N}\left(\textbf{\textit{m}},\textbf{\textit{K}}\right),
$$

- The covariance matrix **K** is constructed based on the chosen inputs  $\{ {\bf x}^{(1)},...,{\bf x}^{(n)} \}$ .
- Entry  $K_{ij}$  is computed by  $k(\mathbf{x}^{(i)}, \mathbf{x}^{(j)})$ .
- Technically, for every choice of inputs  $\{ \mathbf{x}^{(1)},...,\mathbf{x}^{(n)} \}$ , K needs to be positive semi-definite in order to be a valid covariance matrix.
- A function *k*(., .) satisfying this property is called **positive definite**.

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# **COVARIANCE FUNCTION OF A GP / 2**

• Recall, the purpose of the covariance function is to control to which degree the following is fulfilled:

> If two points  $\mathbf{x}^{(i)}$ ,  $\mathbf{x}^{(j)}$  are close in  $\mathcal{X}$ -space, their function values *f*(**x** (*i*) ), *f*(**x** (*j*) ) should be close (**correlated**!) in  $\mathcal Y$ -space.

Closeness of two points  $\mathbf{x}^{(i)}$ ,  $\mathbf{x}^{(j)}$  in input space  $\mathcal{X}$  is measured in terms of  $\boldsymbol{d} = \mathbf{x}^{(i)} - \mathbf{x}^{(j)}$ :

$$
k(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = k(\mathbf{d})
$$



# **COVARIANCE FUNCTION OF A GP: EXAMPLE**

- Let  $f(\mathbf{x})$  be a GP with  $k(\mathbf{x}, \mathbf{x}') = \exp(-\frac{1}{2})$  $\frac{1}{2}$  $||d||^2$ ) with *d* **= <b>x** − **x**<sup>'</sup>.
- Consider two points  $x^{(1)} = 3$  and  $x^{(2)} = 2.5$ .
- If you want to know how correlated their function values are, compute their correlation!



Covariance Function



# **COVARIANCE FUNCTION OF A GP: EXAMPLE**

Assume we observed a value  $y^{(1)} = -0.8$ , the value of  $y^{(2)}$  should be close under the assumption of the above Gaussian process.



# **COVARIANCE FUNCTION OF A GP: EXAMPLE**

- Let us compare another point  $x^{(3)}$  to the point  $x^{(1)}$
- We again compute their correlation
- Their function values are not very much correlated;  $y^{(1)}$  and  $y^{(3)}$ might be far away from each other





# **COVARIANCE FUNCTIONS**

There are three types of commonly used covariance functions:

 $k(.,.)$  is called stationary if it is as a function of  $\boldsymbol{d} = \boldsymbol{x} - \boldsymbol{x}'$ , we write *k*(*d*).

Stationarity is invariance to translations in the input space:

- $k(x, x + d) = k(0, d)$
- $k(.,.)$  is called isotropic if it is a function of  $r = ||\mathbf{x} \mathbf{x}'||$ , we write *k*(*r*). Isotropy is invariance to rotations of the input space and implies stationarity.
- $k(.,.)$  is a dot product covariance function if  $k$  is a function of  $\boldsymbol{x}^T\boldsymbol{x}'$

### **COMMONLY USED COVARIANCE FUNCTIONS**



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 $K_{\nu}(\cdot)$  is the modified Bessel function of the second kind.

### **COMMONLY USED COVARIANCE FUNCTIONS / 2**



- Random functions drawn from Gaussian processes with a Squared Exponential Kernel (left), Polynomial Kernel (middle), and a Matérn Kernel (right,  $\ell = 1$ ).
- The length-scale hyperparameter determines the "wiggliness" of the function.  $\bullet$
- $\bullet$  For Matérn, the  $\nu$  parameter determines how differentiable the process is.

#### **SQUARED EXPONENTIAL COVARIANCE FUNCTION**

The squared exponential function is one of the most commonly used covariance functions.

$$
k(\mathbf{x}, \mathbf{x}') = \exp\bigg(-\frac{\|\mathbf{x} - \mathbf{x}'\|^2}{2\ell^2}\bigg).
$$

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#### **Properties**:

- It depends merely on the distance  $r = ||\mathbf{x} \mathbf{x}'|| \rightarrow$  isotropic and stationary.
- Infinitely differentiable  $\rightarrow$  sometimes deemed unrealistic for modeling most of the physical processes.

#### **CHARACTERISTIC LENGTH-SCALE**

$$
k(\mathbf{x},\mathbf{x}') = \exp\left(-\frac{1}{2\ell^2}\|\mathbf{x} - \mathbf{x}'\|^2\right)
$$

 $\ell$  is called **characteristic length-scale**. Loosely speaking, the characteristic length-scale describes how far you need to move in input space for the function values to become uncorrelated. Higher  $\ell$  induces smoother functions, lower  $\ell$  induces more wiggly functions.



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### **CHARACTERISTIC LENGTH-SCALE / 2**

For  $p > 2$  dimensions, the squared exponential can be parameterized:

$$
k(\mathbf{x}, \mathbf{x}') = \exp\left(-\frac{1}{2} (\mathbf{x} - \mathbf{x}')^\top \mathbf{M} (\mathbf{x} - \mathbf{x}')\right)
$$

Possible choices for the matrix *M* include

$$
M_1 = \ell^{-2}I
$$
  $M_2 = diag(\ell)^{-2}$   $M_3 = \Gamma\Gamma^{T} + diag(\ell)^{-2}$ 

where  $\ell$  is a *p*-vector of positive values and  $\Gamma$  is a  $p \times k$  matrix.

The 2nd (and most important) case can also be written as

$$
k(\mathbf{d}) = \exp\left(-\frac{1}{2}\sum_{j=1}^p \frac{d_j^2}{l_j^2}\right)
$$

# **CHARACTERISTIC LENGTH-SCALE / 3**

What is the benefit of having an individual hyperparameter  $\ell_i$  for each dimension?

- The  $\ell_1, \ldots, \ell_p$  hyperparameters play the role of **characteristic length-scales**.
- Loosely speaking, ℓ*<sup>i</sup>* describes how far you need to move along axis *i* in input space for the function values to be uncorrelated.
- Such a covariance function implements **automatic relevance determination** (ARD), since the inverse of the length-scale  $\ell_i$ determines the relevancy of input feature *i* to the regression.
- If  $\ell_i$  is very large, the covariance will become almost independent of that input, effectively removing it from inference.
- If the features are on different scales, the data can be automatically **rescaled** by estimating  $\ell_1, \ldots, \ell_p$

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### **CHARACTERISTIC LENGTH-SCALE / 4**



For the first plot, we have chosen  $M = I$ : the function varies the same in all directions. The second plot is for  $\textit{M} = \text{diag}(\ell)^{-2}$  and  $\ell = (1,3)$ : The function varies less rapidly as a function of  $x_2$  than  $x_1$  as the length-scale for  $x_1$  is less. In the third plot  $\pmb{M} = \mathsf{\Gamma} \mathsf{\Gamma}^{\mathsf{T}} + \mathsf{diag}(\pmb{\ell})^{-2}$  for  $\mathsf{\Gamma} = (1,-1)^{\top}$  and  $\ell = (6,6)^{\top}.$  Here  $\mathsf{\Gamma}$  gives the direction of the most rapid variation. (Image from Rasmussen & Williams, 2006)