

Outline



- Introduction
- Gaussian processes

Learning objectives

At the end of this lecture, you should be able to:

- Understand the differences between the weight-space and function-space viewpoints of regression
- Explain how GPs are a generalization of the multivariate normal to infinitely many variables
- Understand the different properties of covariance functions and their usefulness
- Perform exact inference in GPs for regression with a Gaussian likelihood
- Estimate the hyper-parameters of a GP (e.g., covariance function and likelihood parameters)
- Explain how GPs can be generalized to work with other likelihood functions (e.g., for classification) and the implications for Bayesian inference

Regression

- Consider models of the inputs $\mathbf{x} \in \mathbb{R}^D$ for continuous response variables $y \in \mathbb{R}$ of the form

$$y = f(\mathbf{x}) + \epsilon, \quad \epsilon \sim \mathcal{N}(\epsilon|0, \sigma^2)$$

- Previously, we assumed f to be a **linear parametric** function of the inputs \mathbf{x}

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

- \mathbf{w} was a D -dimensional vector of parameters (one weight per input dimension)

Regression

- Consider models of the inputs $\mathbf{x} \in \mathbb{R}^D$ for continuous response variables $y \in \mathbb{R}$ of the form

$$y = f(\mathbf{x}) + \epsilon, \quad \epsilon \sim \mathcal{N}(\epsilon|0, \sigma^2)$$

- Previously, we assumed f to be a **linear parametric** function of the inputs \mathbf{x}

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

- \mathbf{w} was a D -dimensional vector of parameters (one weight per input dimension)
- We could therefore write the likelihood for a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$ as

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n|\mathbf{w}^T \mathbf{x}_n, \sigma^2)$$

where $\mathbf{y} = \{y_1, \dots, y_N\}$ and $\mathbf{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$

Frequentist vs Bayesian approach

- Likelihood given by

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n|\mathbf{w}^\top \mathbf{x}_n, \sigma^2)$$

- In a **frequentist approach**, we find the parameters \mathbf{w} that maximize the (log) likelihood

$$\hat{\mathbf{w}}_{\text{ML}} = \arg \max_{\mathbf{w}} \left(\sum_{n=1}^N \log p(y_n|\mathbf{w}, \mathbf{x}_n) \right)$$

- This is called **maximum likelihood (ML)** estimation

Frequentist vs Bayesian approach

- Likelihood given by

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n|\mathbf{w}^\top \mathbf{x}_n, \sigma^2)$$

- In a **frequentist approach**, we find the parameters \mathbf{w} that maximize the (log) likelihood

$$\hat{\mathbf{w}}_{\text{ML}} = \arg \max_{\mathbf{w}} \left(\sum_{n=1}^N \log p(y_n|\mathbf{w}, \mathbf{x}_n) \right)$$

- This is called **maximum likelihood (ML)** estimation
- We can make predictions for new test inputs \mathbf{x}_* by plugging in the estimate $\hat{\mathbf{w}}_{\text{ML}}$

$$p(y_*|\hat{\mathbf{w}}_{\text{ML}}, \mathbf{x}_*)$$

- Point prediction given by

$$\hat{y}_* = (\hat{\mathbf{w}}_{\text{ML}})^\top \mathbf{x}_*$$

Frequentist vs Bayesian approach

- Likelihood given by

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n | \mathbf{w}^T \mathbf{x}_n, \sigma^2)$$

- We can further consider a prior on \mathbf{w} : $p(\mathbf{w}) = \mathcal{N}(\mathbf{w}|\mathbf{0}, \lambda\mathbf{I})$

$$\hat{\mathbf{w}}_{\text{MAP}} = \arg \max_{\mathbf{w}} \left(\sum_{n=1}^N \log p(y_n | \mathbf{w}, \mathbf{x}_n) + \log p(\mathbf{w}) \right)$$

- This is called **maximum-a-posteriori (MAP)** estimation
- Term $\log p(\mathbf{w})$ acts as a penalty term - **regularization**

Frequentist vs Bayesian approach

- Likelihood given by

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n | \mathbf{w}^T \mathbf{x}_n, \sigma^2)$$

- We can further consider a prior on \mathbf{w} : $p(\mathbf{w}) = \mathcal{N}(\mathbf{w}|\mathbf{0}, \lambda\mathbf{I})$

$$\hat{\mathbf{w}}_{\text{MAP}} = \arg \max_{\mathbf{w}} \left(\sum_{n=1}^N \log p(y_n | \mathbf{w}, \mathbf{x}_n) + \log p(\mathbf{w}) \right)$$

- This is called **maximum-a-posteriori (MAP)** estimation
- Term $\log p(\mathbf{w})$ acts as a penalty term - **regularization**
- As before, we make predictions by plugging in the estimate $\hat{\mathbf{w}}_{\text{MAP}}$

$$p(y_* | \hat{\mathbf{w}}_{\text{MAP}}, \mathbf{x}_*)$$

- Point prediction given by

$$\hat{y}_* = (\hat{\mathbf{w}}_{\text{MAP}})^T \mathbf{x}_*$$

Frequentist vs Bayesian approach

- Likelihood given by

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n|\mathbf{w}^T \mathbf{x}_n, \sigma^2)$$

- Prior given by $p(\mathbf{w}) = \mathcal{N}(\mathbf{w}|\mathbf{0}, \lambda\mathbf{I})$
- In a **Bayesian approach**, we treat \mathbf{w} as a latent variable and do inference on it

$$p(\mathbf{w}|\mathbf{y}, \mathbf{X}) = \frac{p(\mathbf{y}|\mathbf{w}, \mathbf{X}) p(\mathbf{w})}{p(\mathbf{y}|\mathbf{X})}$$

- We obtain a **full posterior distribution** on \mathbf{w} rather than a point estimate!

Frequentist vs Bayesian approach

- Likelihood given by

$$p(\mathbf{y}|\mathbf{w}, \mathbf{X}) = \prod_{n=1}^N \mathcal{N}(y_n|\mathbf{w}^T \mathbf{x}_n, \sigma^2)$$

- Prior given by $p(\mathbf{w}) = \mathcal{N}(\mathbf{w}|\mathbf{0}, \lambda\mathbf{I})$
- In a **Bayesian approach**, we treat \mathbf{w} as a latent variable and do inference on it

$$p(\mathbf{w}|\mathbf{y}, \mathbf{X}) = \frac{p(\mathbf{y}|\mathbf{w}, \mathbf{X}) p(\mathbf{w})}{p(\mathbf{y}|\mathbf{X})}$$

- We obtain a **full posterior distribution** on \mathbf{w} rather than a point estimate!
- We can make predictions for new test input \mathbf{x}_* by **averaging** over the values of \mathbf{w}

$$p(y_*|\mathbf{x}_*, \mathbf{y}, \mathbf{X}) = \int p(y_*|\mathbf{w}, \mathbf{x}_*) p(\mathbf{w}|\mathbf{y}, \mathbf{X}) d\mathbf{w}$$

- Marginal likelihood given by

$$p(\mathbf{y}|\mathbf{X}) = \int p(\mathbf{y}|\mathbf{w}, \mathbf{X}) p(\mathbf{w}) d\mathbf{w}$$

Weight-space vs. function-space view

- Consider a dataset of target variables $\mathbf{y} = \{y_1, \dots, y_N\}$ and their corresponding inputs $\mathbf{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$
- In Bayesian linear regression, we assumed that

$$y = f(\mathbf{x}) + \epsilon, \quad \epsilon \sim \mathcal{N}(0, \sigma^2)$$

where $f(\mathbf{x}) = \mathbf{w}^\top \mathbf{x}$

- We placed a prior on the weights $p(\mathbf{w})$ and performed inference to compute its posterior distribution $p(\mathbf{w}|\mathbf{y}, \mathbf{X})$

Weight-space vs. function-space view

- Consider a dataset of target variables $\mathbf{y} = \{y_1, \dots, y_N\}$ and their corresponding inputs $\mathbf{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$
- In Bayesian linear regression, we assumed that

$$y = f(\mathbf{x}) + \epsilon, \quad \epsilon \sim \mathcal{N}(0, \sigma^2)$$

where $f(\mathbf{x}) = \mathbf{w}^\top \mathbf{x}$

- We placed a prior on the weights $p(\mathbf{w})$ and performed inference to compute its posterior distribution $p(\mathbf{w}|\mathbf{y}, \mathbf{X})$
- We can consider this in vector-form

$$\mathbf{y} = \mathbf{f} + \epsilon, \quad \epsilon \sim \mathcal{N}(\mathbf{0}, \sigma^2 \mathbf{I})$$

where $\mathbf{f} = f(\mathbf{X}) = \mathbf{w}^\top \mathbf{X}$

Weight-space vs. function-space view

- Consider a dataset of target variables $\mathbf{y} = \{y_1, \dots, y_N\}$ and their corresponding inputs $\mathbf{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$
- In Bayesian linear regression, we assumed that

$$y = f(\mathbf{x}) + \epsilon, \quad \epsilon \sim \mathcal{N}(0, \sigma^2)$$

where $f(\mathbf{x}) = \mathbf{w}^\top \mathbf{x}$

- We placed a prior on the weights $p(\mathbf{w})$ and performed inference to compute its posterior distribution $p(\mathbf{w}|\mathbf{y}, \mathbf{X})$
- We can consider this in vector-form

$$\mathbf{y} = \mathbf{f} + \epsilon, \quad \epsilon \sim \mathcal{N}(\mathbf{0}, \sigma^2 \mathbf{I})$$

where $\mathbf{f} = f(\mathbf{X}) = \mathbf{w}^\top \mathbf{X}$

- Can we avoid \mathbf{w} altogether and model $p(\mathbf{f})$ directly?
- Instead of working with weights \mathbf{w} , can we work with the functions $f(\mathbf{x})$?
I.e. put a prior on \mathbf{f} and perform inference on it?

Playtime!

- Jupyter notebook: "12 - Gaussian processes.ipynb"
- Part 1: From multivariate Gaussians to Gaussian processes

From multivariate Gaussians to Gaussian processes

- **Definition:** a Gaussian process (GP) is a collection of random variables, any finite number of which have (consistent) joint Gaussian distributions

From multivariate Gaussians to Gaussian processes

- **Definition:** a Gaussian process (GP) is a collection of random variables, any finite number of which have (consistent) joint Gaussian distributions
- Consider a model of the form

$$y = f(\mathbf{x}) + \epsilon$$

- Now consider a multivariate (joint) Gaussian distribution over the N-dimensional vector $\mathbf{f} = (f(\mathbf{x}_1), \dots, f(\mathbf{x}_N))^T$

$$\mathbf{f} \sim \mathcal{N}(\mathbf{f}|\boldsymbol{\mu}, \boldsymbol{\Sigma})$$

- A multivariate Gaussian distribution is fully specified by a mean vector $\boldsymbol{\mu}$ and a covariance matrix $\boldsymbol{\Sigma}$

From multivariate Gaussians to Gaussian processes

- **Definition:** a Gaussian process (GP) is a collection of random variables, any finite number of which have (consistent) joint Gaussian distributions
- Consider a model of the form

$$y = f(\mathbf{x}) + \epsilon$$

- Now consider a multivariate (joint) Gaussian distribution over the N-dimensional vector $\mathbf{f} = (f(\mathbf{x}_1), \dots, f(\mathbf{x}_N))^T$

$$\mathbf{f} \sim \mathcal{N}(\mathbf{f}|\boldsymbol{\mu}, \boldsymbol{\Sigma})$$

- A multivariate Gaussian distribution is fully specified by a mean vector $\boldsymbol{\mu}$ and a covariance matrix $\boldsymbol{\Sigma}$
- A GP is a **stochastic process** fully specified by a mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ and a positive definite covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]$
- Therefore, a GP is a **generalization** of a multivariate Gaussian distribution to infinitely many variables

Gaussian processes

- A GP is a **stochastic process** fully specified by a mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ and a positive definite covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]$
- Mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ determines the mean of any arbitrary point \mathbf{x} in the input space
 - Commonly assumed to be a zero-value vector, i.e. $m(\mathbf{x}) = 0$

Gaussian processes

- A GP is a **stochastic process** fully specified by a mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ and a positive definite covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]]$
- Mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ determines the mean of any arbitrary point \mathbf{x} in the input space
 - Commonly assumed to be a zero-value vector, i.e. $m(\mathbf{x}) = 0$
- Covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]]$ determines how any two points in the input space covary (often called *kernels*)
 - Specifies basic aspects of the process such as smoothness, periodicity, stationarity and isotropy

Gaussian processes

- A GP is a **stochastic process** fully specified by a mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ and a positive definite covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]$
- Mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ determines the mean of any arbitrary point \mathbf{x} in the input space
 - Commonly assumed to be a zero-value vector, i.e. $m(\mathbf{x}) = 0$
- Covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]$ determines how any two points in the input space covary (often called *kernels*)
 - Specifies basic aspects of the process such as smoothness, periodicity, stationarity and isotropy
- If we loosely see a function as a infinitely long vector \mathbf{f} , then we can think of a GP as a **probability distribution over functions!**

Gaussian processes

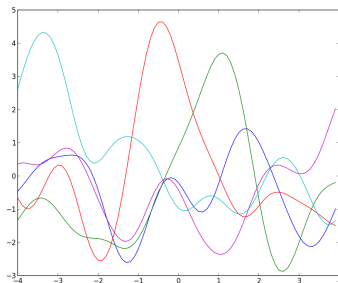
- A GP is a **stochastic process** fully specified by a mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ and a positive definite covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]$
- Mean function $m(\mathbf{x}) = \mathbb{E}[f(\mathbf{x})]$ determines the mean of any arbitrary point \mathbf{x} in the input space
 - Commonly assumed to be a zero-value vector, i.e. $m(\mathbf{x}) = 0$
- Covariance function $k(\mathbf{x}, \mathbf{x}') = \text{cov}[f(\mathbf{x}), f(\mathbf{x}')]$ determines how any two points in the input space covary (often called *kernels*)
 - Specifies basic aspects of the process such as smoothness, periodicity, stationarity and isotropy
- If we loosely see a function as a infinitely long vector \mathbf{f} , then we can think of a GP as a **probability distribution over functions!**
- **Main idea:** we place a GP prior over the function values \mathbf{f} ; together with some likelihood function, we compute the GP posterior (we will return to this later...)
- GPs are **Bayesian non-parametric** models!

Covariance functions

- Most common choice is the **squared exponential (SE)**

$$k_{\text{SE}}(\mathbf{x}, \mathbf{x}') = \exp \left(- \sum_{d=1}^D \frac{(x_d - x'_d)^2}{2l^2} \right)$$

- Also called Gaussian kernel, RBF kernel, exponentiated quadratic, etc.
- Parameter l defining the characteristic length-scale
- Goes to unity as \mathbf{x} becomes closer to \mathbf{x}'
- Nearby points are more likely to covary!
- GP prior with a SE covariance function prefers **smooth functions**



Covariance functions

- Other popular covariance functions:
 - Periodic (PER) covariance function

$$k_{\text{PER}}(\mathbf{x}, \mathbf{x}') = h^2 \exp \left(- \frac{1}{2\ell^2} \sin^2 \left(\frac{\pi}{p} \sum_{d=1}^D (x_d - x'_d) \right) \right)$$

where h controls the amplitude and p is the period

Covariance functions

- Other popular covariance functions:
 - Periodic (PER) covariance function

$$k_{\text{PER}}(\mathbf{x}, \mathbf{x}') = h^2 \exp \left(- \frac{1}{2\ell^2} \sin^2 \left(\frac{\pi}{p} \sum_{d=1}^D (x_d - x'_d) \right) \right)$$

where h controls the amplitude and p is the period

- White noise (WN) covariance function (with variance σ^2)

$$k_{\text{WN}}(\mathbf{x}, \mathbf{x}') = \sigma^2 \delta(\mathbf{x}, \mathbf{x}')$$

where $\delta(\mathbf{x}, \mathbf{x}')$ is the Kronecker delta function (1 when $\mathbf{x} = \mathbf{x}'$, 0 otherwise)

Covariance functions

- Other popular covariance functions:
 - Periodic (PER) covariance function

$$k_{\text{PER}}(\mathbf{x}, \mathbf{x}') = h^2 \exp \left(- \frac{1}{2\ell^2} \sin^2 \left(\frac{\pi}{p} \sum_{d=1}^D (x_d - x'_d) \right) \right)$$

where h controls the amplitude and p is the period

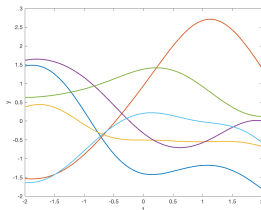
- White noise (WN) covariance function (with variance σ^2)

$$k_{\text{WN}}(\mathbf{x}, \mathbf{x}') = \sigma^2 \delta(\mathbf{x}, \mathbf{x}')$$

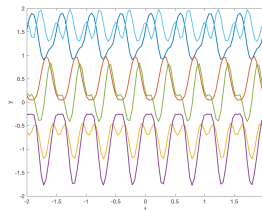
where $\delta(\mathbf{x}, \mathbf{x}')$ is the Kronecker delta function (1 when $\mathbf{x} = \mathbf{x}'$, 0 otherwise)

- Sums and products of proper covariance function are also valid covariance functions!

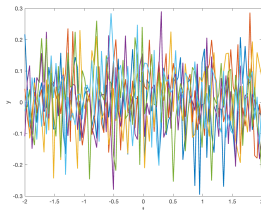
Covariance functions



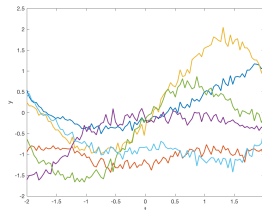
(a) squared exponential (SE)



(b) periodic (PER)



(c) white noise (WN)



(d) SE+WN

Figure: Samples from Gaussian processes with different covariance functions.

Constructing a GP

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$
- Build covariance matrix \mathbf{K} , where $\mathbf{K}_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$

Constructing a GP

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$
- Build covariance matrix \mathbf{K} , where $\mathbf{K}_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$
- Specifies a multivariate Gaussian distribution on \mathbf{f}

$$\mathbf{f} \sim \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

- This is our prior distribution over \mathbf{f}
- We can use it to sample from the GP prior!

Playtime!

- Jupyter notebook: "12 - Gaussian processes.ipynb"
- Part 2: Sampling from a GP with different covariance functions

Inference with a GP

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$
- Build covariance matrix \mathbf{K} , where $\mathbf{K}_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$
- Specifies a multivariate Gaussian distribution on \mathbf{f}

$$\mathbf{f} \sim \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

- This is our prior distribution over \mathbf{f} , $p(\mathbf{f})$, but we **need a likelihood too!**

Inference with a GP

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$
- Build covariance matrix \mathbf{K} , where $\mathbf{K}_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$
- Specifies a multivariate Gaussian distribution on \mathbf{f}

$$\mathbf{f} \sim \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

- This is our prior distribution over \mathbf{f} , $p(\mathbf{f})$, but we **need a likelihood too!**
- For continuous outputs $y \in \mathbb{R}$, an obvious choice is a Gaussian likelihood:

$$p(y_n|f_n) = \mathcal{N}(y_n|f_n, \sigma^2)$$

Inference with a GP

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$
- Build covariance matrix \mathbf{K} , where $\mathbf{K}_{ij} = k(\mathbf{x}_i, \mathbf{x}_j)$
- Specifies a multivariate Gaussian distribution on \mathbf{f}

$$\mathbf{f} \sim \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

- This is our prior distribution over \mathbf{f} , $p(\mathbf{f})$, but we **need a likelihood too!**
- For continuous outputs $y \in \mathbb{R}$, an obvious choice is a Gaussian likelihood:

$$p(y_n | f_n) = \mathcal{N}(y_n | f_n, \sigma^2)$$

- Using Bayes rule, we can compute the posterior over \mathbf{f} (exact inference)

$$p(\mathbf{f}|\mathbf{y}, \mathbf{X}) = \frac{p(\mathbf{y}|\mathbf{f}) p(\mathbf{f})}{p(\mathbf{y}|\mathbf{X})}$$

- Compare this equation with the posterior for \mathbf{w} in Bayesian linear regression

Marginal likelihood

- Making use of marginalization property for Gaussian distributions (see slide 10 of lecture 9), the marginal distribution of \mathbf{y} is given by

$$\begin{aligned} p(\mathbf{y}|\mathbf{X}) &= \int \underbrace{p(\mathbf{y}|\mathbf{f})}_{\text{likelihood}} \underbrace{p(\mathbf{f}|\mathbf{X})}_{\text{GP prior}} d\mathbf{f} \\ &= \int \mathcal{N}(\mathbf{y}|\mathbf{f}, \sigma^2 \mathbf{I}) \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K}) d\mathbf{f} \\ &= \mathcal{N}(\mathbf{y}|\mathbf{0}, \sigma^2 \mathbf{I} + \mathbf{K}) \end{aligned}$$

Marginal likelihood

- Making use of marginalization property for Gaussian distributions (see slide 10 of lecture 9), the marginal distribution of \mathbf{y} is given by

$$\begin{aligned} p(\mathbf{y}|\mathbf{X}) &= \int \underbrace{p(\mathbf{y}|\mathbf{f})}_{\text{likelihood}} \underbrace{p(\mathbf{f}|\mathbf{X})}_{\text{GP prior}} d\mathbf{f} \\ &= \int \mathcal{N}(\mathbf{y}|\mathbf{f}, \sigma^2\mathbf{I}) \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K}) d\mathbf{f} \\ &= \mathcal{N}(\mathbf{y}|\mathbf{0}, \sigma^2\mathbf{I} + \mathbf{K}) \end{aligned}$$

- We can use $p(\mathbf{y}|\mathbf{X})$ to optimize the parameters θ of the covariance function!

$$\hat{\theta} = \arg \max_{\theta} \left(\log \mathcal{N}(\mathbf{y}|\mathbf{0}, \sigma^2\mathbf{I} + \mathbf{K}) \right)$$

Note

This another example of maximum marginal likelihood (also called type-II maximum likelihood, or empirical Bayes) from lecture 11.

Making predictions

- Our aim is to make a prediction y_* for a new input \mathbf{x}_*

Making predictions

- Our aim is to make a prediction y_* for a new input \mathbf{x}_*
- The joint distribution over y_*, y_1, \dots, y_N is simply given by

$$p(y_*, \mathbf{y} | \mathbf{x}_*, \mathbf{X}) = \mathcal{N}(y_*, \mathbf{y} | \mathbf{0}, \mathbf{V})$$

with

$$\mathbf{V} = \begin{pmatrix} \sigma^2 \mathbf{I} + \mathbf{K} & \mathbf{k}_* \\ \mathbf{k}_*^\top & \sigma^2 + k_{**} \end{pmatrix}$$

where $\mathbf{k}_* = k(\mathbf{x}, x_*)$ and $k_{**} = k(x_*, x_*)$

Making predictions

- Our aim is to make a prediction y_* for a new input \mathbf{x}_*
- The joint distribution over y_*, y_1, \dots, y_N is simply given by

$$p(y_*, \mathbf{y} | \mathbf{x}_*, \mathbf{X}) = \mathcal{N}(y_*, \mathbf{y} | \mathbf{0}, \mathbf{V})$$

with

$$\mathbf{V} = \begin{pmatrix} \sigma^2 \mathbf{I} + \mathbf{K} & \mathbf{k}_* \\ \mathbf{k}_*^\top & \sigma^2 + k_{**} \end{pmatrix}$$

where $\mathbf{k}_* = k(\mathbf{x}, x_*)$ and $k_{**} = k(x_*, x_*)$

- From the joint distribution, we can now determine the distribution of y_* , i.e. the predictive distribution:

$$p(y_* | \mathbf{y}, \mathbf{x}_*, \mathbf{X}) = \mathcal{N}(y_* | \mathbf{k}_*^\top (\sigma^2 \mathbf{I} + \mathbf{K})^{-1} \mathbf{y}, k_{**} + \sigma^2 - \mathbf{k}_*^\top (\sigma^2 \mathbf{I} + \mathbf{K})^{-1} \mathbf{k}_*)$$

- Note: this is a direct application of the conditional probability for Gaussians (see slide 11 from lecture 9)

Playtime!

- Jupyter notebook: "12 - Gaussian processes.ipynb"
- Part 3: Inference and maximum marginal likelihood optimization

GP classification

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- What if y_n is discrete?

GP classification

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- What if y_n is discrete?
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$, such that

$$p(\mathbf{f}) = \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

GP classification

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- What if y_n is discrete?
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$, such that

$$p(\mathbf{f}) = \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

- For binary outputs $y_n \in \{0, 1\}$, a possible choice is the Probit function $\Phi(f_n)$:

$$p(y_n|f_n) = \Phi(f_n) = \int_{-\infty}^{f_n} \mathcal{N}(u|0, 1) du$$

GP classification

- Given a dataset $\mathcal{D} = \{\mathbf{x}_n, y_n\}_{n=1}^N$
- What if y_n is discrete?
- Define GP prior for function values \mathbf{f} : $\mathbf{f} \sim \mathcal{GP}(m(\mathbf{x}) = 0, k(\mathbf{x}, \mathbf{x}'))$, such that

$$p(\mathbf{f}) = \mathcal{N}(\mathbf{f}|\mathbf{0}, \mathbf{K})$$

- For binary outputs $y_n \in \{0, 1\}$, a possible choice is the Probit function $\Phi(f_n)$:

$$p(y_n|f_n) = \Phi(f_n) = \int_{-\infty}^{f_n} \mathcal{N}(u|0, 1) du$$

- Exact inference is no longer tractable; must resort to approximate methods

$$p(\mathbf{f}|\mathbf{y}, \mathbf{X}) = \frac{(\prod_{n=1}^N \Phi(f_n)) p(\mathbf{f})}{p(\mathbf{y}|\mathbf{X})}$$

- Similar approaches can be used to handle other types of outputs
 - Real, binary, categorical, positive real, positive integer or ordinal responses

Learning more about GPs

- A Visual Exploration of Gaussian Processes - **This notebook is a must!**
<https://distill.pub/2019/visual-exploration-gaussian-processes/>
- Videolecture: Gaussian Processes, C. Rasmussen.
http://videlectures.net/mlss09uk_rasmussen_gp/
- Book: Gaussian Processes for Machine Learning, C. Rasmussen and C. Williams.
Free! <http://www.gaussianprocess.org/gpml/>
- Book: Pattern Recognition and Machine Learning, C. Bishop.