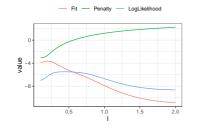
# **Introduction to Machine Learning**

# **Gaussian Processes Training of a Gaussian Process**





#### Learning goals

- Training of GPs via Maximum Likelihood estimation of its hyperparameters
- Computational complexity is governed by matrix inversion of the covariance matrix

## TRAINING OF A GAUSSIAN PROCESS

- To make predictions for a regression task by a Gaussian process, one simply needs to perform matrix computations.
- But for this to work out, we assume that the covariance functions is fully given, including all of its hyperparameters.
- A very nice property of GPs is that we can learn the numerical hyperparameters of a selected covariance function directly during GP training.



# TRAINING A GP VIA MAXIMUM LIKELIHOOD

Let us assume

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

where  $f(\mathbf{x}) \sim \mathcal{GP}(\mathbf{0}, k(\mathbf{x}, \mathbf{x}'|\boldsymbol{\theta}))$ .

Observing  ${\it y}\sim \mathcal{N}\left({\it 0},{\it K}+\sigma^2{\it I}\right)$ , the marginal log-likelihood (or evidence) is

$$\log p(\mathbf{y} \mid \mathbf{X}, \boldsymbol{\theta}) = \log \left[ (2\pi)^{-n/2} |\mathbf{K}_{y}|^{-1/2} \exp \left( -\frac{1}{2} \mathbf{y}^{\top} \mathbf{K}_{y}^{-1} \mathbf{y} \right) \right]$$
$$= -\frac{1}{2} \mathbf{y}^{T} \mathbf{K}_{y}^{-1} \mathbf{y} - \frac{1}{2} \log |\mathbf{K}_{y}| - \frac{n}{2} \log 2\pi.$$

with  $K_y := K + \sigma^2 I$  and  $\theta$  denoting the hyperparameters (the parameters of the covariance function).



## TRAINING A GP VIA MAXIMUM LIKELIHOOD / 2

The three terms of the marginal likelihood have interpretable roles, considering that the model becomes less flexible as the length-scale increases:

- the data fit  $-\frac{1}{2} \mathbf{y}^T \mathbf{K}_y^{-1} \mathbf{y}$ , which tends to decrease if the length scale increases
- the complexity penalty  $-\frac{1}{2}\log |\mathbf{K}_y|$ , which depends on the covariance function only and which increases with the length-scale, because the model gets less complex with growing length-scale
- a normalization constant  $-\frac{n}{2} \log 2\pi$



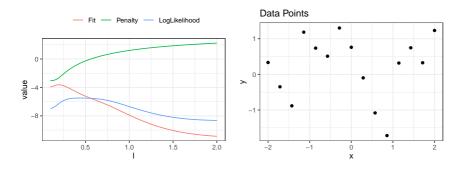
To visualize this, we consider a zero-mean Gaussian process with squared exponential kernel

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

- $\bullet$  Recall, the model is smoother and less complex for higher length-scale  $\ell.$
- We show how the
  - data fit  $-\frac{1}{2} \mathbf{y}^T \mathbf{K}_{y}^{-1} \mathbf{y}$ ,
  - the complexity penalty  $-\frac{1}{2} \log |K_y|$ , and
  - ullet the overall value of the marginal likelihood  $\log p(m{y} \mid m{X}, m{ heta})$

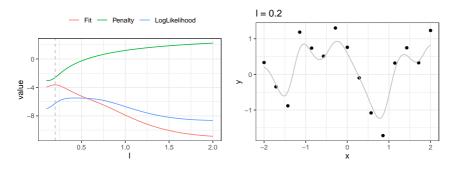
behave for increasing value of  $\ell$ .







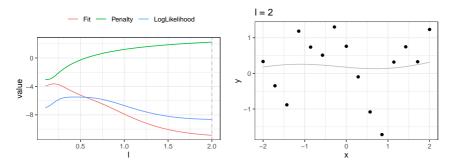
The left plot shows how values of the data fit  $-\frac{1}{2} \mathbf{y}^T \mathbf{K}_y^{-1} \mathbf{y}$ , the complexity penalty  $-\frac{1}{2} \log |\mathbf{K}_y|$  (high value means less penalization) and the overall marginal likelihood  $\log p(\mathbf{y} \mid \mathbf{X}, \boldsymbol{\theta})$  behave for increasing values of  $\ell$ .





The left plot shows how values of the data fit  $-\frac{1}{2} \pmb{y}^T \pmb{K}_y^{-1} \pmb{y}$ , the complexity penalty  $-\frac{1}{2} \log |\pmb{K}_y|$  (high value means less penalization) and the overall marginal likelihood  $\log p(\pmb{y} \mid \pmb{X}, \pmb{\theta})$  behave for increasing values of  $\ell$ .

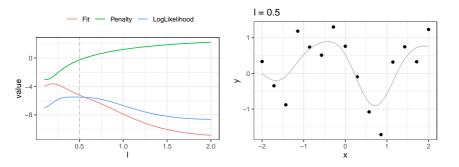
A small  $\ell$  results in a good fit, but a high complexity penalty (low  $-\frac{1}{2} \log |\mathbf{K}_y|$ ).





The left plot shows how values of the data fit  $-\frac{1}{2} \mathbf{y}^T \mathbf{K}_y^{-1} \mathbf{y}$ , the complexity penalty  $-\frac{1}{2} \log |\mathbf{K}_y|$  (high value means less penalization) and the overall marginal likelihood  $\log p(\mathbf{y} \mid \mathbf{X}, \boldsymbol{\theta})$  behave for increasing values of  $\ell$ .

A large  $\ell$  results in a poor fit.





The left plot shows how values of the data fit  $-\frac{1}{2} \mathbf{y}^T \mathbf{K}_y^{-1} \mathbf{y}$ , the complexity penalty  $-\frac{1}{2} \log |\mathbf{K}_y|$  (high value means less penalization) and the overall marginal likelihood  $\log p(\mathbf{y} \mid \mathbf{X}, \boldsymbol{\theta})$  behave for increasing values of  $\ell$ .

The maximizer of the log-likelihood,  $\ell=0.5$ , balances complexity and fit.

#### TRAINING A GP VIA MAXIMUM LIKELIHOOD

using  $\frac{\partial}{\partial \theta_i} \mathbf{K}^{-1} = -\mathbf{K}^{-1} \frac{\partial \mathbf{K}}{\partial \theta_i} \mathbf{K}^{-1}$  and  $\frac{\partial}{\partial \theta} \log |\mathbf{K}| = \operatorname{tr} \left( \mathbf{K}^{-1} \frac{\partial \mathbf{K}}{\partial \theta} \right)$ .

To set the hyperparameters by maximizing the marginal likelihood, we seek the partial derivatives w.r.t. the hyperparameters

$$\frac{\partial}{\partial \theta_{j}} \log p(\mathbf{y} \mid \mathbf{X}, \boldsymbol{\theta}) = \frac{\partial}{\partial \theta_{j}} \left( -\frac{1}{2} \mathbf{y}^{T} \mathbf{K}_{y}^{-1} \mathbf{y} - \frac{1}{2} \log |\mathbf{K}_{y}| - \frac{n}{2} \log 2\pi \right) 
= \frac{1}{2} \mathbf{y}^{T} \mathbf{K}^{-1} \frac{\partial \mathbf{K}}{\partial \theta_{j}} \mathbf{K}^{-1} \mathbf{y} - \frac{1}{2} \operatorname{tr} \left( \mathbf{K}^{-1} \frac{\partial \mathbf{K}}{\partial \boldsymbol{\theta}} \right) 
= \frac{1}{2} \operatorname{tr} \left( (\mathbf{K}^{-1} \mathbf{y} \mathbf{y}^{T} \mathbf{K}^{-1} - \mathbf{K}^{-1}) \frac{\partial \mathbf{K}}{\partial \theta_{j}} \right)$$



## TRAINING A GP VIA MAXIMUM LIKELIHOOD / 2

- The complexity and the runtime of training a Gaussian process is dominated by the computational task of inverting K - or let's rather say for decomposing it.
- Standard methods require  $\mathcal{O}(n^3)$  time (!) for this.
- Once  $K^{-1}$  or rather the decomposition -is known, the computation of the partial derivatives requires only  $\mathcal{O}(n^2)$  time per hyperparameter.
- Thus, the computational overhead of computing derivatives is small, so using a gradient based optimizer is advantageous.



## TRAINING A GP VIA MAXIMUM LIKELIHOOD / 3

Workarounds to make GP estimation feasible for big data include:

- using kernels that yield sparse *K*: cheaper to invert.
- subsampling the data to estimate  $\theta$ :  $\mathcal{O}(m^3)$  for subset of size m.
- combining estimates on different subsets of size m: Bayesian committee,  $\mathcal{O}(nm^2)$ .
- using low-rank approximations of K by using only a representative subset ("inducing points") of m training data  $X_m$ :

  Nyström approximation  $K \approx K_{nm}K_{mm}^-K_{mn}$ ,  $\mathcal{O}(nmk + m^3)$  for a rank-k-approximate inverse of  $K_{mm}$ .
- exploiting structure in K induced by the kernel: exact solutions but complicated maths, not applicable for all kernels.

... this is still an active area of research.

