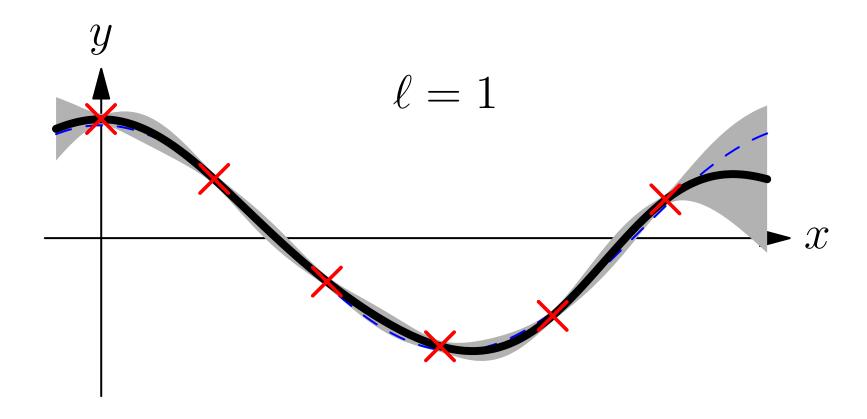
# **Advanced Machine Learning**

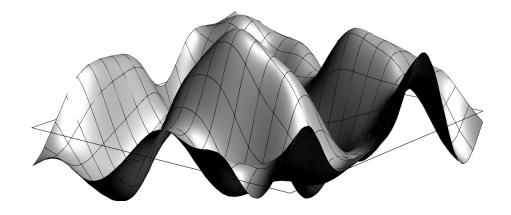
### Gaussian Processes



 $Gaussian\ Processes,\ regression$ 

### **Outline**

- 1. Introduction
- 2. Gaussian Processes
- 3. Bayesian Inference
- 4. Hyper-parameters



- Gaussian processes (GPs) are a mathematically defined ensemble of functions
- They can be combined with Bayesian inference to give one of the most powerful regression techniques
- Although Bayesian they can be used in a black-box fashion due to the ubiquity of the prior
- Mathematically they are a bit complicated

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- (You can use Gaussian Processes for classification, e.g. by inferring the probabilities of being in a class, but we ignore this as regression is where GP excel)
- In regression we have some p dimensional feature vectors  $m{x}_i$  and some target  $y_i \in \mathbb{R}$
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- We can think of a solution as a function f(x)
- We can put a prior probability distribution, p(f), on a function, f, that prefers smooth functions
- We can then compute a posterior probability distribution on functions given the data,  $p(f|\mathcal{D})$
- As a likelihood,  $p(y_i|f(x_i))$ , we use the probability of observing  $y_i$  given the true function value is  $f(x_i)$
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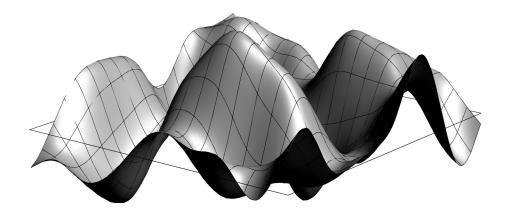
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- 2. Gaussian Processes
- 3. Bayesian Inference
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- Gaussian Processes are probability distributions over functions
- (Functions can be viewed as vectors in an infinite dimensional vector space)
- In the Gaussian Process,  $\mathcal{GP}(m,k)$ , the probability of a function, f, is proportional

$$p(f|m,k) \propto e^{-\frac{1}{2} \int (f(\boldsymbol{x}) - m(\boldsymbol{x})) k^{-1}(\boldsymbol{x},\boldsymbol{y}) (f(\boldsymbol{y}) - m(\boldsymbol{y})) d\boldsymbol{x} d\boldsymbol{y}}$$

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## Meaning of GP

- ullet To understand GP's we can discretise space,  $oldsymbol{x}$ , into a lattice of points  $\{oldsymbol{x}_i\}$
- Then (assuming  $m(\boldsymbol{x}) = 0$ )

$$p(f|m,k) \propto \prod_{i} e^{-\frac{f_i^2 k^{-1}(\boldsymbol{x}_i, \boldsymbol{x}_i)}{2}} + f_i \sum_{j} k^{-1}(\boldsymbol{x}_i, \boldsymbol{x}_j) f_j$$

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$$\mathbb{E}\left[\left(f(\boldsymbol{x}) - m(\boldsymbol{x})\right)\left(f(\boldsymbol{y}) - m(\boldsymbol{y})\right)\right] = k(\boldsymbol{x}, \boldsymbol{y})$$

- This is sometimes know as a kernel—it must be positive semi-definite
- It is a free "parameter" that the user gets to choose (although we can learn its parameters too)
- If  $k(\boldsymbol{x}, \boldsymbol{y})$  is a function of  $\boldsymbol{x} \boldsymbol{y}$  it is "stationary"
- If  $k(\boldsymbol{x}, \boldsymbol{y})$  is a function of  $\|\boldsymbol{x} \boldsymbol{y}\|$  it is also "isometric"

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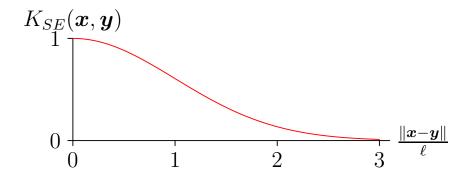
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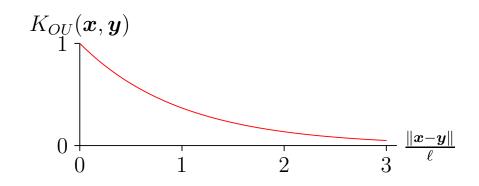
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# Popular Choices of GP Kernel Function

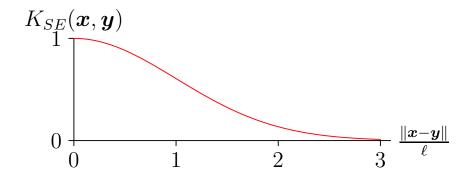
- Constant:  $k_{\rm C}(\boldsymbol{x},\boldsymbol{y}) = C$
- Gaussian noise:  $k_{\rm GN}({\boldsymbol x},{\boldsymbol y}) = \sigma^2 \delta_{{\boldsymbol x},{\boldsymbol y}}$
- Squared exponential:  $k_{\mathrm{SE}}(m{x},m{y}) = \exp\left(-\frac{\|m{x}-m{y}\|^2}{2\ell^2}\right)$
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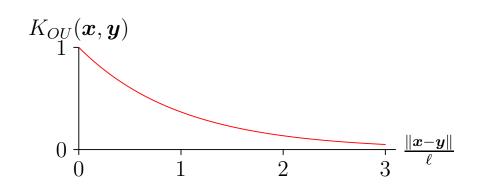




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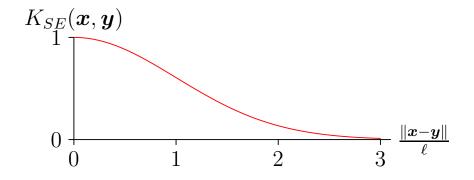
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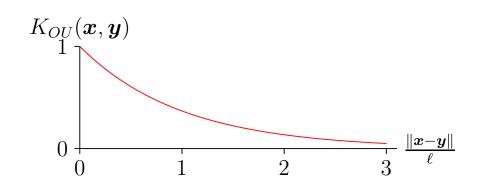




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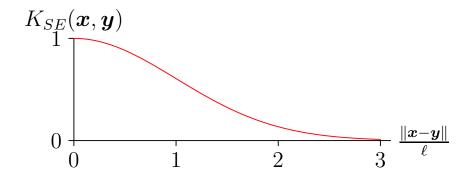
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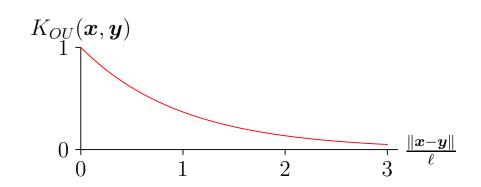




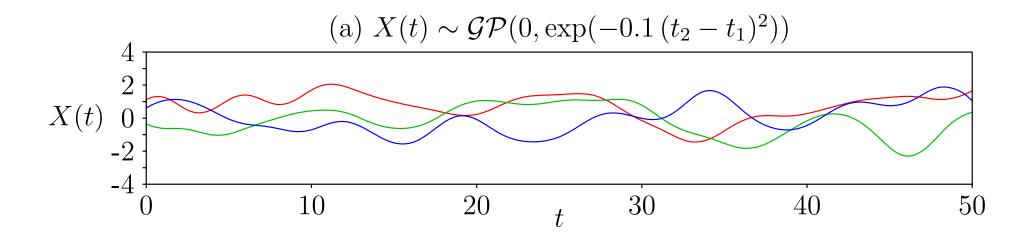
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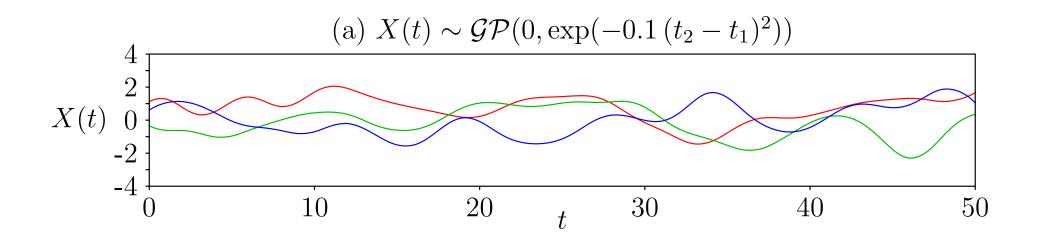


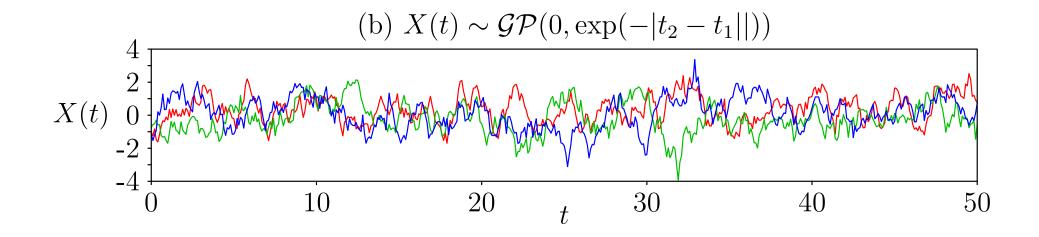


#### **Gaussian Process Worlds**

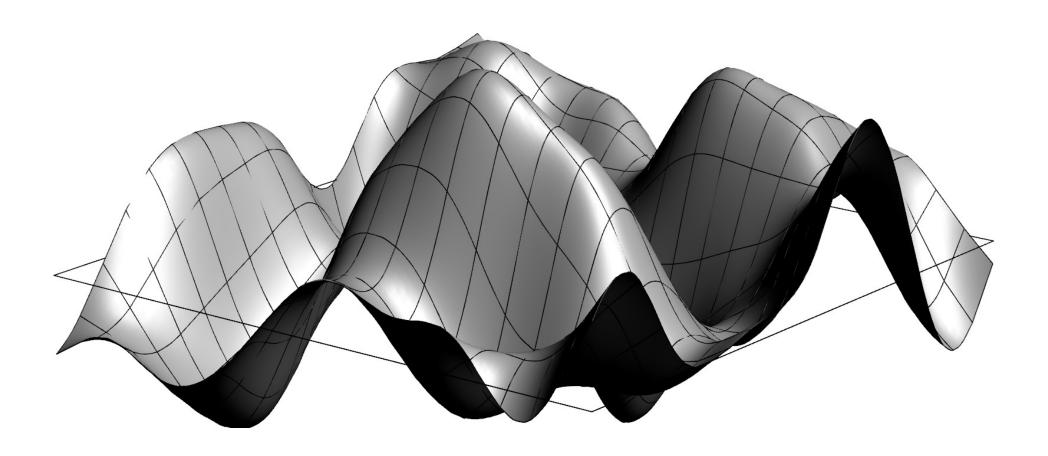


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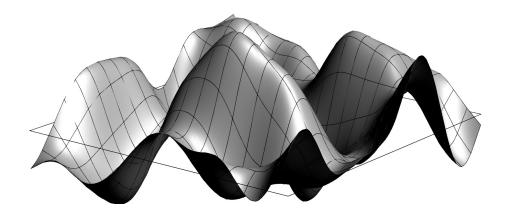


### 2-D Gaussian Processes



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- 2. Gaussian Processes
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• Given some data points  $\mathcal{D} = ((\boldsymbol{x}_i, y_i) | i = 1, ..., m)$  the likelihood (assuming Gaussian error are independence of the data point) is given by

$$p(\mathcal{D}|f) = \prod_{i=1}^{m} \mathcal{N}(y_i | f(\boldsymbol{x}_i), \sigma^2)$$

- Using a Gausssian Process prior we can compute a posterior using Bayes's rule
- The posterior is a Gaussian Process with a shifted mean and variance depending on the data-points
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- Denoting the matrices of covariances between data points as  ${\bf K}$  with elements  $k({m x}_i,{m x}_j)$
- Denoting the covariance between the data points and a particular position,  $x_*$  as  $k_*$  with elements  $k(x_i,x_*)$
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$$p(\boldsymbol{y}, f_*) = \mathcal{N}\left(\begin{pmatrix} \boldsymbol{y} \\ f_* \end{pmatrix} \middle| \boldsymbol{0}, \begin{pmatrix} \mathbf{K} + \sigma^2 \mathbf{I} & \boldsymbol{k}_* \\ \boldsymbol{k}_*^\mathsf{T} & k_* \end{pmatrix}\right)$$

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- ullet Then the distribution of function values at points at  $oldsymbol{x}_i$  and  $oldsymbol{x}_*$  is

$$p(\boldsymbol{y}, f_*) = \mathcal{N}\left(\begin{pmatrix} \boldsymbol{y} \\ f_* \end{pmatrix} \middle| \boldsymbol{0}, \begin{pmatrix} \mathbf{K} + \sigma^2 \mathbf{I} & \boldsymbol{k}_* \\ \boldsymbol{k}_*^\mathsf{T} & k_* \end{pmatrix}\right)$$

• To compute the posterior  $p(f_*|y)$  we use

$$p(f_*|\mathbf{y}) = \frac{p(f_*,\mathbf{y})}{p(\mathbf{y})}$$

- where  $p(\boldsymbol{y}) = \int p(f_*, \boldsymbol{y}) df_*$
- Because all integrals are Gaussian we can compute the integral to obtain

$$p(f_*|\boldsymbol{y}) = \mathcal{N}\left(f_* \middle| \boldsymbol{k}_*^\mathsf{T} (\mathbf{K} + \sigma^2 \mathbf{I})^{-1} \boldsymbol{y}, k - \boldsymbol{k}_*^\mathsf{T} (\mathbf{K} + \sigma^2 \mathbf{I})^{-1} \boldsymbol{k}_*\right)$$

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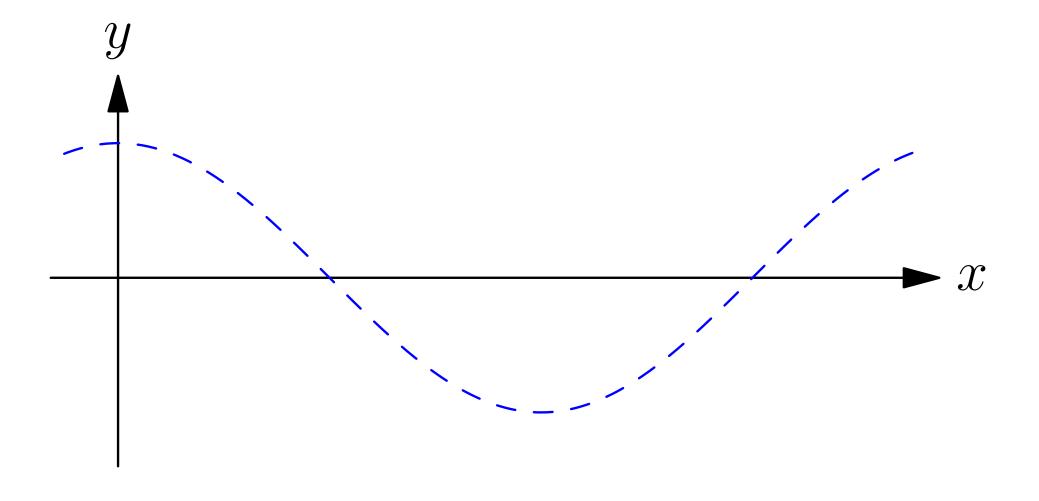
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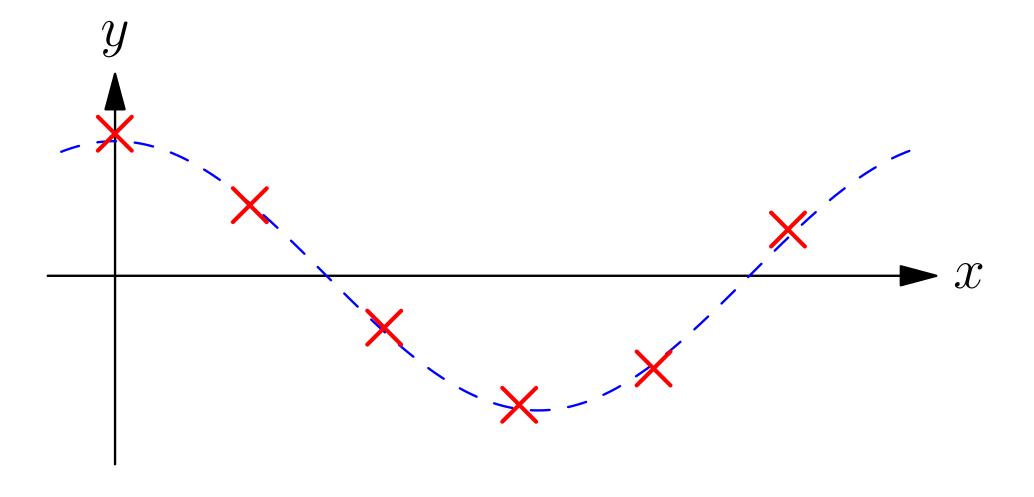
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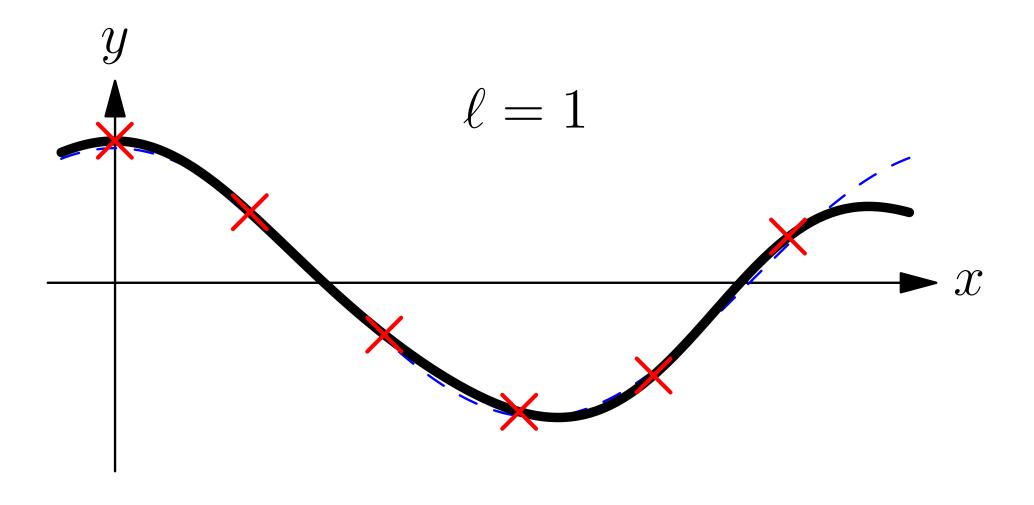
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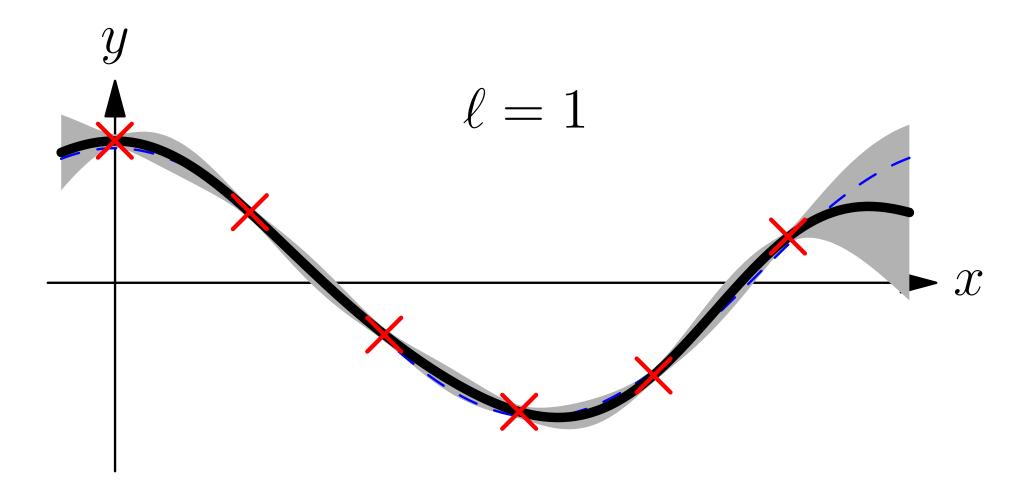
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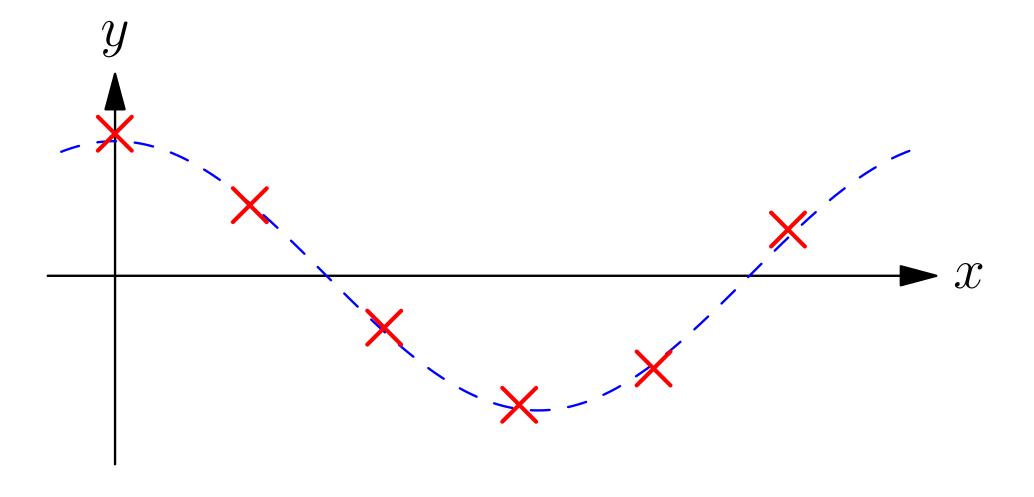
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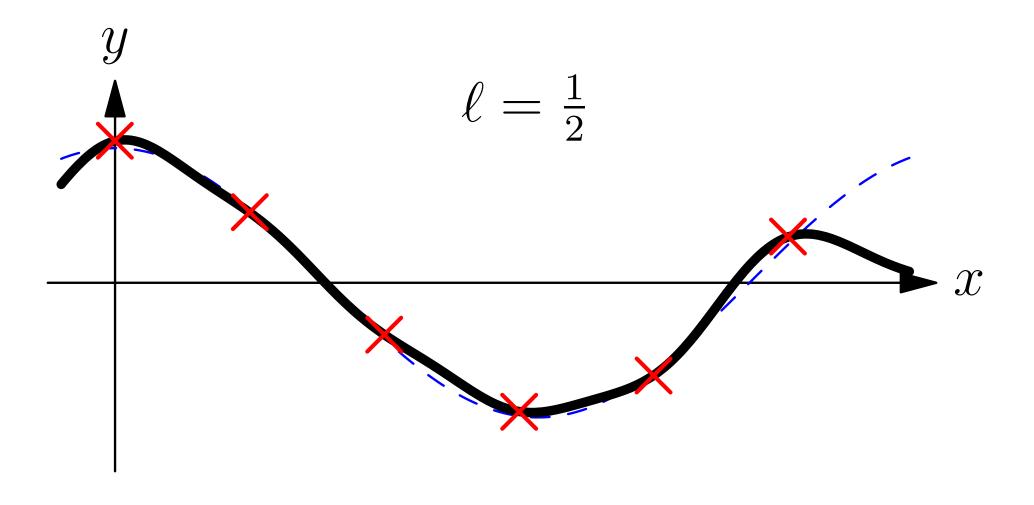
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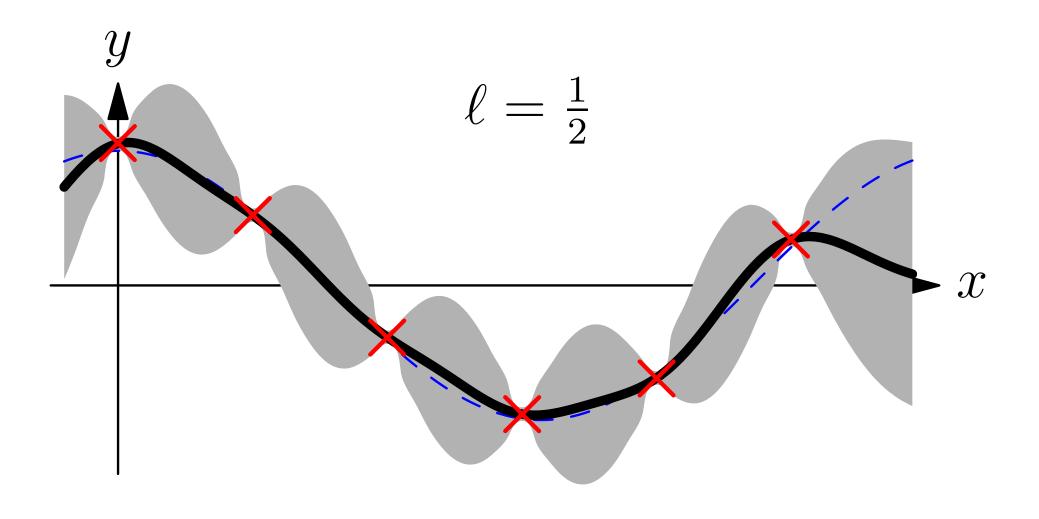
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- This might be used with a time series
- ullet The much more typical situation in machine learning is for x to have many features so we are doing multi-dimensional regression
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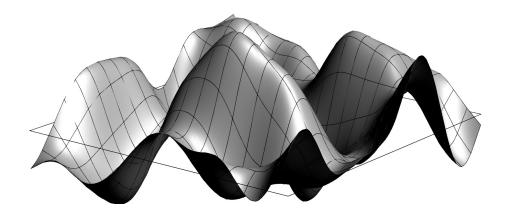
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#### **Outline**

- 1. Introduction
- 2. Gaussian Processes
- 3. Bayesian Inference
- 4. Hyper-parameters



- Choosing the correct covariance function is critical
- Most covariance functions include a continuous **hyper-parameter** (e.g. the correlation length  $\ell$ ) that we have to choose correctly
- This is typical of many Bayesian problems were we have some set of hyper-parameters,  $\phi$ , describing the model
- These are different to the normal parameters we learn (e.g. weights  ${m w}$  or in GP the functions  $f({m x})$ )
- In Bayesian inference we learn the posterior for these normal parameters

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#### **Evidence Framework**

• The normalisation factor,  $p(\mathcal{D}|\phi)$  is known as the **marginal** likelihood or evidence

$$p(\mathcal{D}|\boldsymbol{\phi}) = \int p(\mathcal{D}|f,\boldsymbol{\phi})p(f|\boldsymbol{\phi})df$$

ullet We can perform a Bayesian calculation at a second level by putting a prior on  $\phi$ 

$$p(\phi|\mathcal{D}) = \frac{p(\mathcal{D}|\phi)p(\phi)}{p(\mathcal{D})}$$

From this we can now marginalise out the hyper-parameters

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- \* Second term measure complexity of model
- ★ Last term is a common normalisation constant
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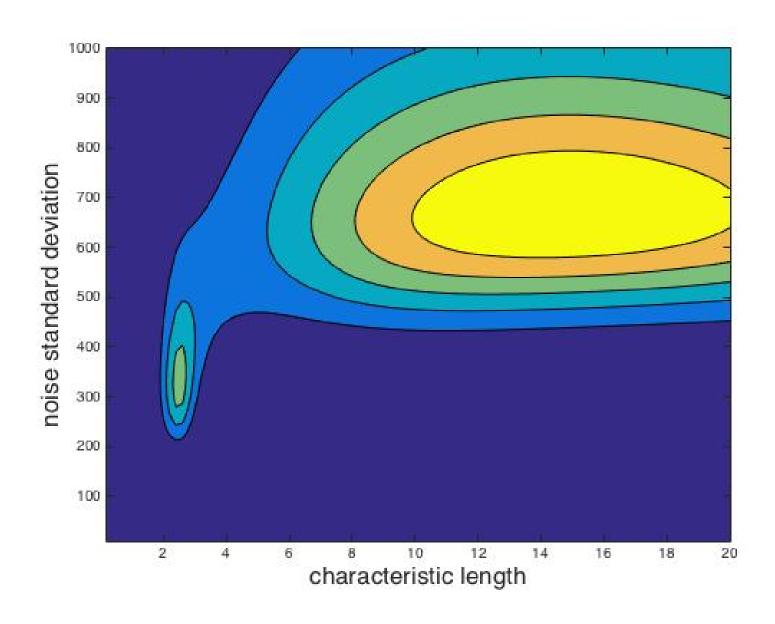
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- Could overfit!

# Example (slightly pathological)



- Gaussian processes are very powerful for regression (and classification?)
- Because all calculations involve Gaussian integrals we can compute everything in closed form
- (Actually its a pain to do the mathematics because you end up working with inverse of matrices)
- Fairly generic (black-box) technique because the prior captures many continuity constraints
- We can use the evidence framework (probability of data) to do model selection and hyper-parameter optimisations

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