

# Gaussian Process

## Gaussian / Normal Distribution

- Univariate Form

$$X \sim N(\mu, \sigma^2)$$
$$p(x) = \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right)$$

- Multivariate Form

$$X \sim N(X|\mu, \Sigma)$$
$$X \in \mathbb{R}^d \quad \mu \in \mathbb{R}^d \quad \Sigma \in \mathbb{R}^{d \times d} \quad \Sigma_{ij} = E(x_i - \mu_i)(x_j - \mu_j) = \text{cov}(x_i, x_j)$$
$$p(X) = \left(\frac{1}{\sqrt{2\pi}}\right)^d \cdot \frac{1}{|\Sigma|^{\frac{1}{2}}} \cdot \exp\left(-\frac{1}{2}(X - \mu)^T \Sigma^{-1}(X - \mu)\right)$$

- $\Sigma$  is a covariance matrix, determines how  $x_i, x_j$  increase together or not
- CLT
  - When  $n$  independent random variables are summed up, their normalized sum tends a *Gaussian distributed random variable*, even if these original random variables are not Gaussian
- Any **linear combination** of Gaussian distributed random variables follow a Gaussian distribution
- **Concatenation** of Gaussian distributed random variables result in a multivariate Gaussian distributed random variable
- Given  $x = \begin{pmatrix} x_a \\ x_b \end{pmatrix} \in \mathbb{R}^{a+b}$ , if

$$x \sim N\left(\begin{pmatrix} \mu_a \\ \mu_b \end{pmatrix}, \begin{pmatrix} \Sigma_{aa} & \Sigma_{ab} \\ \Sigma_{ba} & \Sigma_{bb} \end{pmatrix}\right)$$

then

- i.  $x_a \sim N(\mu_a, \Sigma_{aa}), x_b \sim N(\mu_b, \Sigma_{bb})$
- ii.  $P(x_a|x_b) = N(x_a|\mu_{a|b}, \Sigma_{a|b})$   
 $\mu_{a|b} = \mu_a + \Sigma_{ab}\Sigma_{bb}^{-1}(x_b - \mu_b)$   
 $\Sigma_{a|b} = \Sigma_{aa} - \Sigma_{ab}\Sigma_{bb}^{-1}\Sigma_{ba}$

## Different views of Linear Regression

### ERM view

- the original form:

$$\min_w \sum_{i=1}^n (y_i - w^T \phi(x_i))^2$$

### MLE view

- $y = w^T \phi(x) + \epsilon, \quad \epsilon \sim N(0, \sigma^2)$ 
  - $\epsilon$  is a Gaussian noise
- $P(y|x; w, \sigma^2) = N(y|w^T \phi(x), \sigma^2)$ 
  - $w$  is the parameter
  - $\sigma^2$  is a hyperparameter
- MLE form:

$$\begin{aligned}
& \max_w \sum_{i=1}^n \log P(y_i|x_i) \\
& \Leftrightarrow \max_w \sum_{i=1}^n \log \left( \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{(y_i - w^T \phi(x_i))^2}{2\sigma^2}\right) \right) \\
& \Leftrightarrow \max_w \sum_{i=1}^n -\frac{1}{2} \log 2\pi\sigma^2 - \frac{(y_i - w^T \phi(x_i))^2}{2\sigma^2} \\
& \Leftrightarrow \min_w \sum_{i=1}^n (y_i - w^T \phi(x_i))^2
\end{aligned}$$

## MAP view

- **Maximum a Posteriori**

- Previously, we treat  $w$  as fixed(constant) parameter. In Bayesian viewpoint, the world is uncertain, even the parameter  $w$  are random variables
- Select the mode of the posteriori distribution as a point estimation

- Prior probability:

$$P(w) = N(w|0, \sigma_w^2 I) \quad w \in \mathbb{R}^d$$

- $\mu = 0$  is a **prior belief**
- 我们先验地认为,  $w$  的期望为0, 方差越小, 说明我们对 $w$ 的取值在0附近这一事件越自信

- Estimation for *Ridge Regression*:

$$P(y|x, w; \sigma^2, \sigma_w^2) = N(y|w^T \phi(x), \sigma^2)$$

From Bayesian Expectation Equation:

$$P(w|y, x) = \frac{P(y|x, w)P(w|x)}{P(y|x)}$$

$P(y|x)$  is not dependent on  $w$ , let it be  $z$ ,  $P(w|x)$  is actually  $P(w)$ , so

$$\begin{aligned}
P(w|y, x) &= \frac{1}{z} \cdot \prod_{i=1}^n P(y_i|x_i, w)P(w) \\
&= \frac{1}{z} \cdot \left(\frac{1}{\sqrt{2\pi}\sigma}\right)^n \cdot \exp\left(-\frac{\sum_{i=1}^n (y_i - w^T \phi(x_i))^2}{2\sigma^2}\right) \cdot \left(\frac{1}{\sqrt{2\pi}\sigma_w}\right)^d \cdot \exp\left(-\frac{w^T w}{2\sigma_w^2}\right) \\
\max_w P(w|y, x) &\Leftrightarrow \max_w -\frac{1}{2\sigma^2} \sum_{i=1}^n (y_i - w^T \phi(x_i))^2 - \frac{1}{2\sigma_w^2} w^T w \\
&\Leftrightarrow \min_w \sum_{i=1}^n (y_i - w^T \phi(x_i))^2 + \frac{\sigma^2}{\sigma_w^2} \|w\|^2
\end{aligned}$$

## Stochastic (Random) Process

- A collection of (infinitely many) random variables along on index set( $\mathbb{N}$  or  $\mathbb{R}$  or  $\mathbb{R}^d \dots$ )
- Infinitely many  $\{(x_i, y_i)\}$  specify a distribution and function  $y(x)$ . Each sample of  $\{(x_i, y_i)\}$  forms a *deterministic function*  $y(x)$
- We can specify a **Random Process** by specifying the **joint distribution** of all random variables

## Gaussian Process (GP)

- We specify the joint distribution over **any finite collection** of variables and require the joint distribution to be Gaussian
- $\{x_1, \dots, x_n\}$  is **any** set of  $n$  points in the index set with sampled value  $\{y_1, \dots, y_n\}$ , then

$$GP \Leftrightarrow y_1, \dots, y_n \text{ have a Multivariate Gaussian Distribution}$$

- We need to specify  $\mu_i = \text{mean}(x_i)$  and  $\Sigma_{ij} = k(x_i, x_j)$  to determine a GP

$$y(x) \sim GP(\text{mean}(x), k(\cdot, \cdot))$$

- we usually use  $\text{mean}(x) = 0$  (prior belief), which means that  $y = w^T \phi(x) = 0$  as  $w \sim N(0, \sigma_w^2 I)$
- we only need  $k(x_i, x_j) = \exp(-\frac{\|x_i - x_j\|^2}{2\sigma^2})$ 
  - $k(x_i, x_j)$  越大, 表明  $x_i$  和  $x_j$  越接近, 就越有可能同增减
- Suppose  $n$  training points  $\{x_1, \dots, x_n\}$ , let  $K$  be Gram matrix,  $K_{ij} = k(x_i, x_j)$ 
  - $k$  is a valid kernel only if  $K \succeq 0$  for any  $\{x_1, \dots, x_n\}$
  - Now given a new test point  $x^*$ , we can compute joint distribution of  $\begin{pmatrix} y^* \\ y \end{pmatrix} \in \mathbb{R}^{n+1}$

$$P\left(\begin{pmatrix} y^* \\ y \end{pmatrix}\right) = N(y|0, \begin{pmatrix} k(x^*, x^*) & k^T(x^*) \\ k(x^*) & K \end{pmatrix}), \quad k(x^*) = \begin{pmatrix} k(x^*, x_1) \\ \vdots \\ k(x^*, x_n) \end{pmatrix}$$

$$P(y^*|y) = N(y^*|\mu^*, \Sigma^*)$$

$$\mu^* = 0 + k^T(x^*)K^{-1}(y - 0) = k^T(x^*)K^{-1}y$$

$$\Sigma^* = k(x^*, x^*) - k^T(x^*)K^{-1}k(x^*)$$

So we have the dual form of Linear Regression:

$$f(x^*) = \mu^* = k^T(x^*)K^{-1}y$$

- $\mu^*$  is called **posterori mean**
- In a more realistic setting, we can only observe  $\hat{y} = y + \epsilon$ ,  $\epsilon \sim N(0, \sigma^2 I)$ , then

$$P(\hat{y}) = N(y|0, K + \sigma^2 I)$$

$$\mu^* = k^T(x^*)(K + \sigma^2 I)^{-1}y$$

$$\Sigma^* = k(x^*, x^*) - k^T(x^*)(K + \sigma^2 I)^{-1}k(x^*)$$

because

$$\begin{aligned} \text{cov}(\hat{y}_i, \hat{y}_j) &= E\hat{y}_i\hat{y}_j - E\hat{y}_iE\hat{y}_j \\ &= E(y_i + \epsilon_i)(y_j + \epsilon_j) \\ &= Ey_iy_j + E\epsilon_i\epsilon_j + Ey_i\epsilon_j + Ey_j\epsilon_i \\ &= Ey_iy_j + E\epsilon_i\epsilon_j \\ &= \text{cov}(y_i, y_j) + \text{cov}(\epsilon_i, \epsilon_j) \\ &= k(x_i, x_j) + \sigma^2 1(i = j) \end{aligned}$$

- 也就是说, 每次采样一些  $x$  (服从高斯分布) 及其对应的  $y$ , 得到一个方程  $y(x)$ , 那么多次采样后, 方程  $y(x)$  也服从高斯分布

## Recall Ridge Regression

- recall the MAP view of Ridge Regression, we have  $y_i = w^T \phi(x_i) + \epsilon_i$ , and vectorize this formula:

$$y = w^T \Phi + \epsilon \quad w \sim N(0, \sigma_w^2 I) \quad \epsilon \sim N(0, \sigma^2 I)$$

For  $y_i$  and  $y_j$ :

$$\begin{aligned} \text{cov}(y_i, y_j) &= Ey_iy_j - Ey_iEy_j \\ &= E(w^T \phi(x_i) + \epsilon_i)(w^T \phi(x_j) + \epsilon_j) \\ &= E(\phi^T(x_i)w w^T \phi(x_j)) + E\epsilon_i\epsilon_j \\ &= \phi^T(x_i)Ew w^T \phi(x_j) + \sigma^2 1(i = j) \\ &= \phi^T(x_i)\text{cov}(w w^T)\phi(x_j) + \sigma^2 1(i = j) \\ &= \sigma_w^2 \phi^T(x_i)\phi(x_j) + \sigma^2 1(i = j) \\ &= \sigma_w^2 k(x_i, x_j) + \sigma^2 1(i = j) \end{aligned}$$

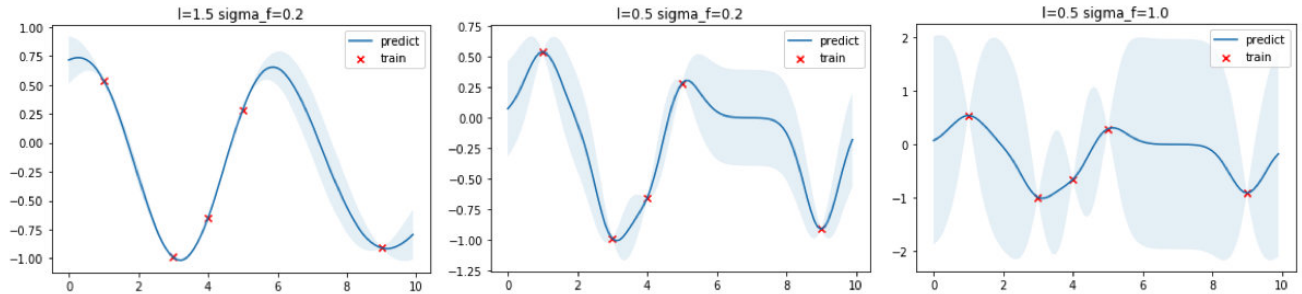
So we have:

$$y \sim N(0, \Sigma) \quad \Sigma = \sigma_w^2 K + \sigma^2 I$$

- For a new test point  $x^* \in \mathbb{R}^{n'}$

$$\begin{aligned}\mu^* &= \sigma_w^2 k^T(x^*)(\sigma_w^2 K + \sigma^2 I)^{-1} y \\ &= k^T(x^*)(K + \frac{\sigma^2}{\sigma_w^2} I)^{-1} y \quad (\text{let } \frac{\sigma^2}{\sigma_w^2} \text{ be } \lambda)\end{aligned}$$

- 贝叶斯视角下的线性回归本质是高斯过程
- GP connects ERM of Ridge Regression, MAP of Linear Regression, and dual solution of Ridge Regression
- **GP for Regression(GPR)**: typically take *RBF* kernel  $k(x, x') = \exp(-\frac{\|x-x'\|^2}{2l^2})$ 
  - $l$  is called *length scale* or *temperature*
  - $\mu(x_i)$  not necessarily equals  $y_i$  (because  $\epsilon_i$  and  $\text{cov}(y_i, y_j)$ )
  - when  $l \rightarrow 0$  and  $\sigma = 0$ ,  $\mu(x_i) = y_i$



## Bayesian Optimization (BO)

- For *Black box* functions (cannot use gradient descent)
  - i. Randomly sample  $n$  points  $x_1, \dots, x_n$  and corresponding  $y_1, \dots, y_n$
  - ii. fit a GP
  - iii. use some acquisition function  $a(x)$  to select next point to evaluate
    - possible  $a(x)$ : *lower confidence bound*  $\mu(x) - \kappa\sigma(x)$
  - iv. Use all  $x, y$  to fit a new GP
  - v. Repeat until reaching a budget
- For hyperparameter optimization, the input is possible value of hyperparameter, and the output is the validation error