

Advanced Machine Learning Lecture 2

Linear Regression

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Announcements

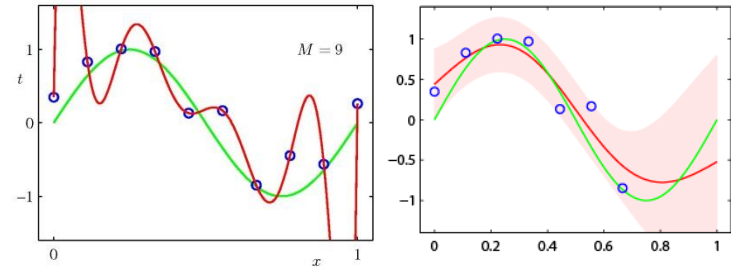
- **No lecture next Monday (22.10.)!**
 - We can use this slot to give a short Matlab tutorial
 - Who would be interested?

This Lecture: *Advanced Machine Learning*

- Regression Approaches

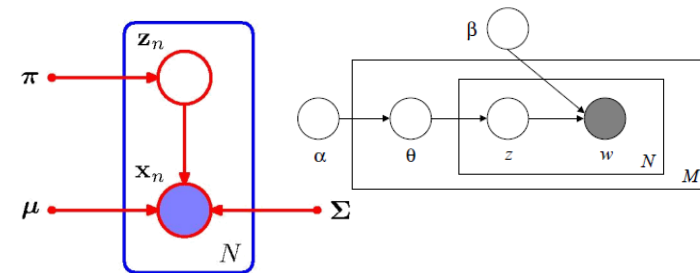
- Linear Regression
- Regularization (Ridge, Lasso)
- Support Vector Regression
- Gaussian Processes

$$f : \mathcal{X} \rightarrow \mathbb{R}$$



- Learning with Latent Variables

- EM and Generalizations
- Dirichlet Processes



- Structured Output Learning

- Large-margin Learning

$$f : \mathcal{X} \rightarrow \mathcal{Y}$$

Topics of This Lecture

- **Recap: Important Concepts from ML Lecture**
 - Probability Theory
 - Bayes Decision Theory
 - Maximum Likelihood Estimation
 - Bayesian Estimation
- **A Probabilistic View on Regression**
 - Least-Squares Estimation as Maximum Likelihood
 - Predictive Distribution
 - Maximum-A-Posteriori (MAP) Estimation
 - Bayesian Curve Fitting
- **Discussion**

Recap: The Rules of Probability

- Basic rules

Sum Rule
$$p(X) = \sum_Y p(X, Y)$$

Product Rule
$$p(X, Y) = p(Y|X)p(X)$$

- From those, we can derive

Bayes' Theorem
$$p(Y|X) = \frac{p(X|Y)p(Y)}{p(X)}$$

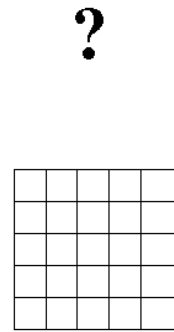
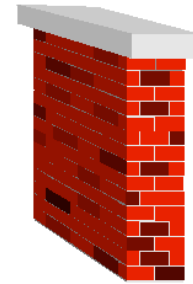
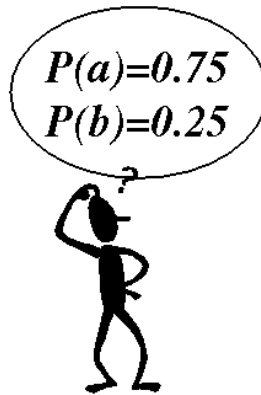
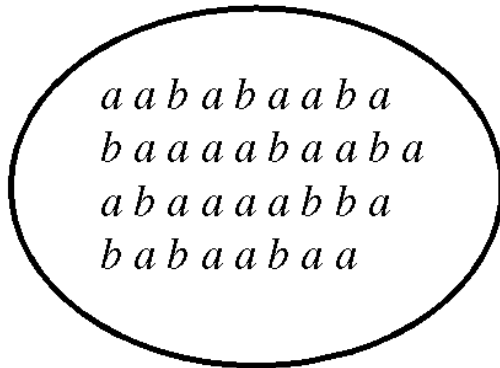
where
$$p(X) = \sum_Y p(X|Y)p(Y)$$

Recap: Bayes Decision Theory

- Concept 1: **Priors** (a priori probabilities)

$$p(C_k)$$

- What we can tell about the probability *before seeing the data*.
- Example:



$$C_1 = a$$

$$p(C_1) = 0.75$$

$$C_2 = b$$

$$p(C_2) = 0.25$$

- In general: $\sum_k p(C_k) = 1$

Recap: Bayes Decision Theory

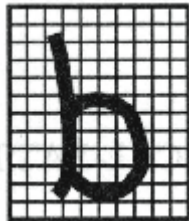
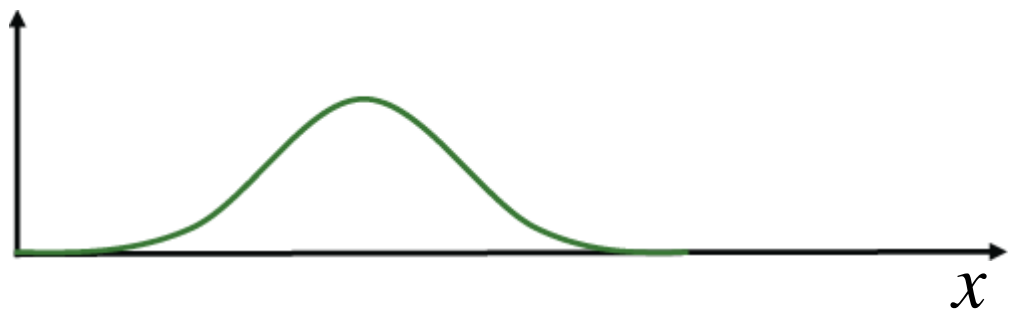
- **Concept 2: Conditional probabilities**

$$p(x | C_k)$$

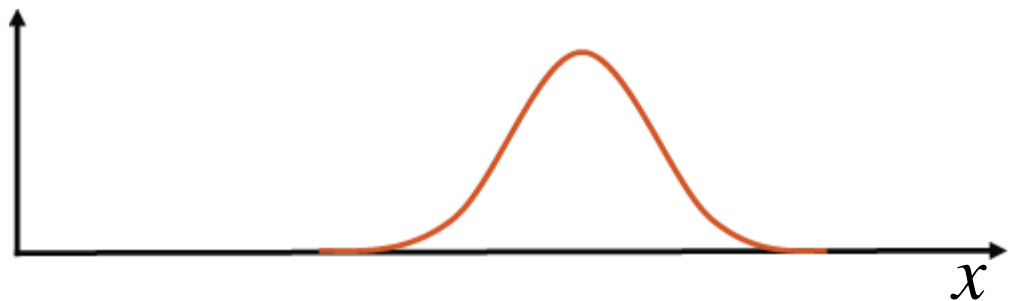
- Let x be a feature vector.
- x measures/describes certain properties of the input.
 - E.g. number of black pixels, aspect ratio, ...
- $p(x|C_k)$ describes its **likelihood** for class C_k .



$$p(x | a)$$



$$p(x | b)$$



Recap: Bayes Decision Theory

- Concept 3: **Posterior probabilities**

$$p(C_k | x)$$

- We are typically interested in the *a posteriori* probability, i.e. the probability of class C_k given the measurement vector x .

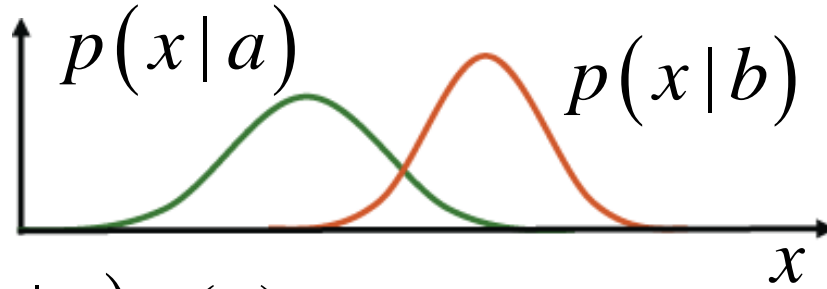
- Bayes' Theorem:

$$p(C_k | x) = \frac{p(x | C_k) p(C_k)}{p(x)} = \frac{p(x | C_k) p(C_k)}{\sum_i p(x | C_i) p(C_i)}$$

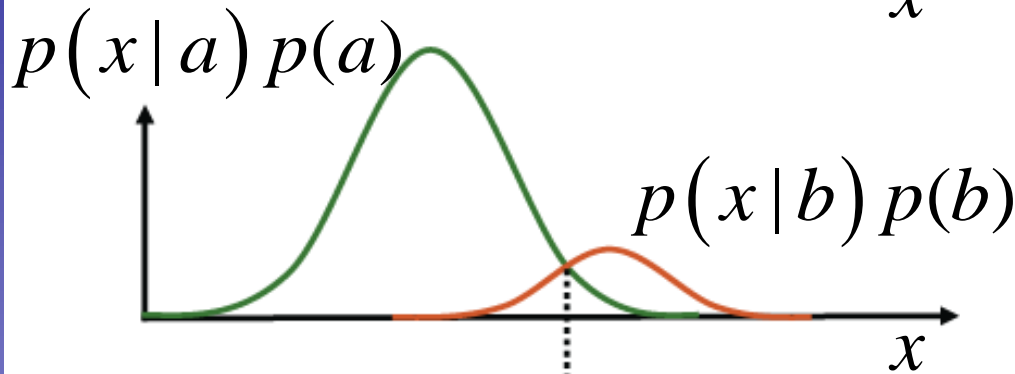
- Interpretation

$$\textit{Posterior} = \frac{\textit{Likelihood} \times \textit{Prior}}{\textit{Normalization Factor}}$$

Recap: Bayes Decision Theory

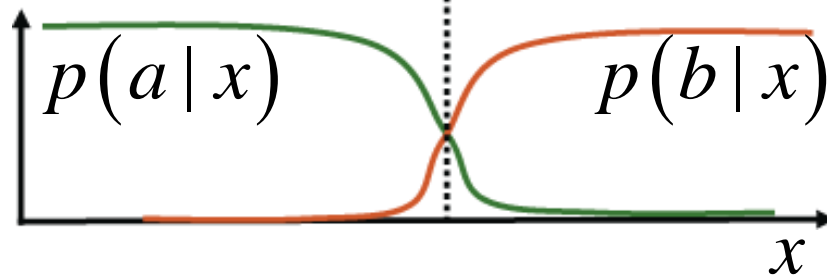


Likelihood



Likelihood \times Prior

Decision boundary



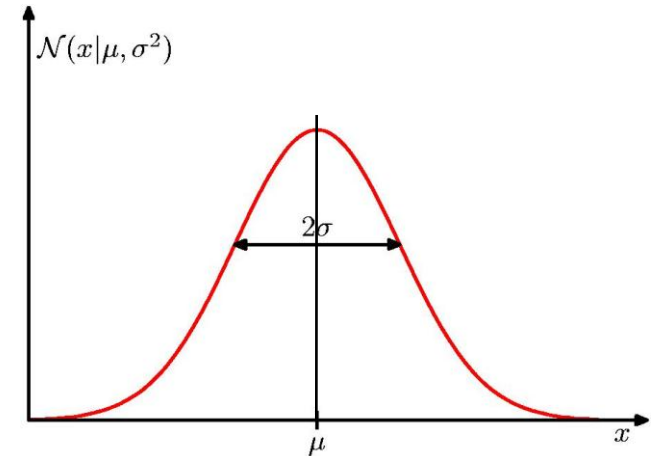
$$\text{Posterior} = \frac{\text{Likelihood} \times \text{Prior}}{\text{NormalizationFactor}}$$

Recap: Gaussian (or Normal) Distribution

- One-dimensional case

- Mean μ
- Variance σ^2

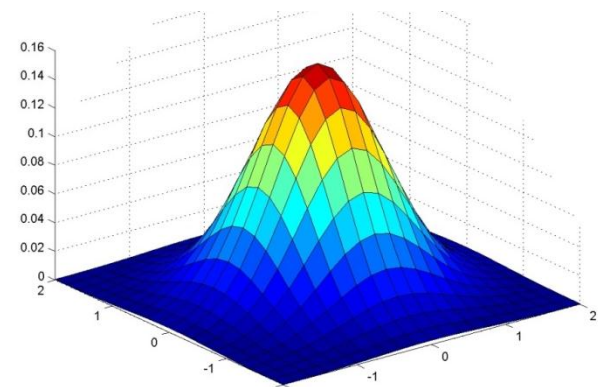
$$\mathcal{N}(x|\mu, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma}} \exp \left\{ -\frac{(x - \mu)^2}{2\sigma^2} \right\}$$



- Multi-dimensional case

- Mean μ
- Covariance Σ

$$\mathcal{N}(\mathbf{x}|\mu, \Sigma) = \frac{1}{(2\pi)^{D/2} |\Sigma|^{1/2}} \exp \left\{ -\frac{1}{2} (\mathbf{x} - \mu)^T \Sigma^{-1} (\mathbf{x} - \mu) \right\}$$



Side Note

- **Notation**

- In many situations, it will be preferable to work with the inverse of the **covariance matrix** Σ :

$$\Lambda = \Sigma^{-1}$$

- We call Λ the **precision matrix**.
- We can therefore also write the Gaussian as

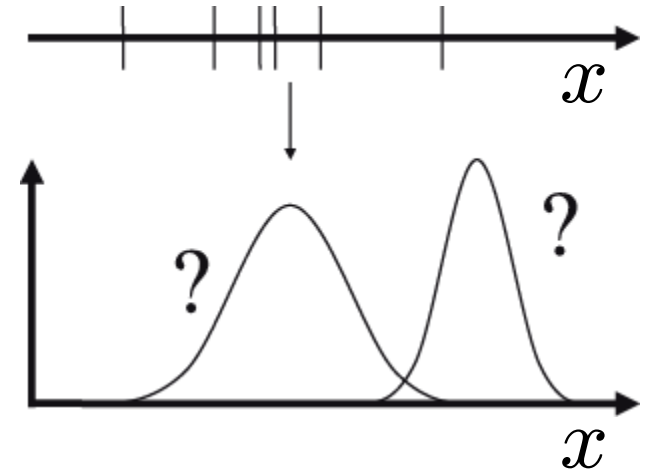
$$\mathcal{N}(x|\mu, \lambda^{-1}) = \frac{1}{\sqrt{2\pi}\lambda^{-1/2}} \exp \left\{ -\frac{\lambda}{2}(x - \mu)^2 \right\}$$

$$\mathcal{N}(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Lambda}^{-1}) = \frac{1}{(2\pi)^{D/2}|\boldsymbol{\Lambda}|^{-1/2}} \exp \left\{ -\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Lambda}(\mathbf{x} - \boldsymbol{\mu}) \right\}$$

Recap: Parametric Methods

- **Given**

- Data $X = \{x_1, x_2, \dots, x_N\}$
- Parametric form of the distribution with parameters θ
- E.g. for Gaussian distrib.: $\theta = (\mu, \sigma)$



- **Learning**

- Estimation of the parameters θ

- **Likelihood of θ**

- Probability that the data X have indeed been generated from a probability density with parameters θ

$$L(\theta) = p(X|\theta)$$

Recap: Maximum Likelihood Approach

- **Computation of the likelihood**

- Single data point: $p(x_n|\theta)$
- Assumption: all data points $X = \{x_1, \dots, x_n\}$ are independent

$$L(\theta) = p(X|\theta) = \prod_{n=1}^N p(x_n|\theta)$$

- **Log-likelihood**

$$E(\theta) = -\ln L(\theta) = -\sum_{n=1}^N \ln p(x_n|\theta)$$

- **Estimation of the parameters θ (Learning)**

- Maximize the likelihood (=minimize the negative log-likelihood)
⇒ Take the derivative and set it to zero.

$$\frac{\partial}{\partial \theta} E(\theta) = -\sum_{n=1}^N \frac{\frac{\partial}{\partial \theta} p(x_n|\theta)}{p(x_n|\theta)} \stackrel{!}{=} 0$$

Recap: Maximum Likelihood Approach

- Applying ML to estimate the parameters of a Gaussian, we obtain

$$\hat{\mu} = \frac{1}{N} \sum_{n=1}^N x_n \quad \text{“sample mean”}$$

- In a similar fashion, we get

$$\hat{\sigma}^2 = \frac{1}{N} \sum_{n=1}^N (x_n - \hat{\mu})^2 \quad \text{“sample variance”}$$

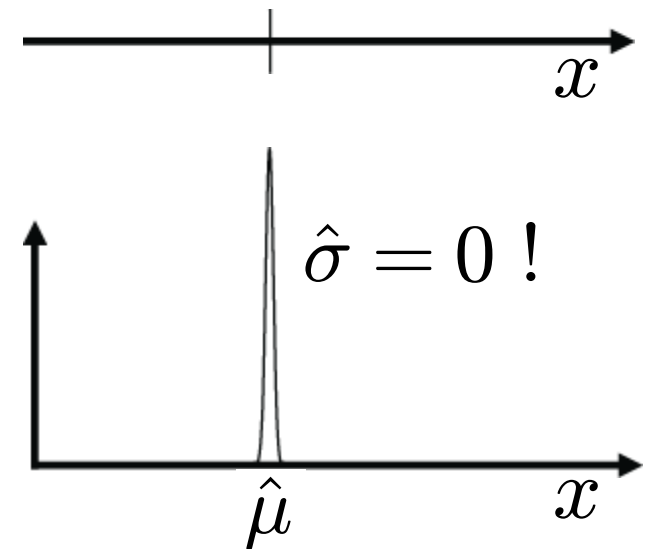
- $\hat{\theta} = (\hat{\mu}, \hat{\sigma})$ is the **Maximum Likelihood estimate** for the parameters of a Gaussian distribution.
- This is a very important result.
- Unfortunately, it is biased...

Recap: Maximum Likelihood - Limitations

- Maximum Likelihood has several significant limitations
 - It systematically underestimates the variance of the distribution!
 - E.g. consider the case

$$N = 1, X = \{x_1\}$$

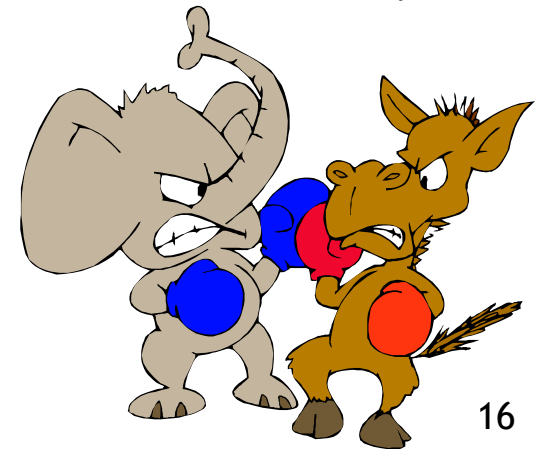
⇒ Maximum-likelihood estimate:



- We say *ML overfits to the observed data*.
- We will still often use ML, but it is important to know about this effect.

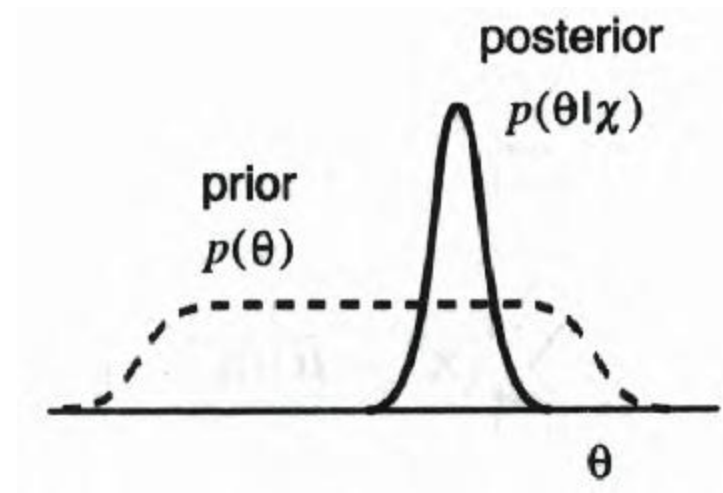
Recap: Deeper Reason

- **Maximum Likelihood** is a **Frequentist** concept
 - In the **Frequentist view**, probabilities are the frequencies of random, repeatable events.
 - These frequencies are fixed, but can be estimated more precisely when more data is available.
- This is in contrast to the **Bayesian** interpretation
 - In the **Bayesian view**, probabilities quantify the uncertainty about certain states or events.
 - This uncertainty can be revised in the light of new evidence.
- **Bayesians and Frequentists do not like each other too well...**



Recap: Bayesian Approach to Learning

- **Conceptual shift**
 - Maximum Likelihood views the true parameter vector θ to be unknown, but fixed.
 - In Bayesian learning, we consider θ to be a random variable.
- This allows us to use knowledge about the parameters θ
 - i.e. to use a prior for θ
 - Training data then converts this prior distribution on θ into a posterior probability density.
- The prior thus encodes knowledge we have about the type of distribution we expect to see for θ .



Recap: Bayesian Learning Approach

- Bayesian view:

- Consider the parameter vector θ as a random variable.
- When estimating the parameters, what we compute is

$$p(x|X) = \int p(x, \theta|X) d\theta$$

Assumption: given θ , this doesn't depend on X anymore

$$p(x, \theta|X) = p(x|\theta, \cancel{X})p(\theta|X)$$

$$p(x|X) = \int \underbrace{p(x|\theta)} p(\theta|X) d\theta$$

This is entirely determined by the parameter θ (i.e. by the parametric form of the pdf).

Recap: Bayesian Learning Approach

$$p(x|X) = \int p(x|\theta)p(\theta|X)d\theta$$

$$p(\theta|X) = \frac{p(X|\theta)p(\theta)}{p(X)} = \frac{p(\theta)}{p(X)}L(\theta)$$

$$p(X) = \int p(X|\theta)p(\theta)d\theta = \int L(\theta)p(\theta)d\theta$$

- Inserting this above, we obtain

$$p(x|X) = \int \frac{p(x|\theta)L(\theta)p(\theta)}{p(X)}d\theta = \int \frac{p(x|\theta)L(\theta)p(\theta)}{\int L(\theta)p(\theta)d\theta}d\theta$$

Recap: Bayesian Learning Approach

- Discussion

Likelihood of the parametric form θ given the data set X .

Estimate for x based on parametric form θ

Prior for the parameters θ

$$p(x|X) = \int \frac{p(x|\theta)L(\theta)p(\theta)}{\int L(\theta)p(\theta)d\theta} d\theta$$

Normalization: integrate over all possible values of θ

- The more uncertain we are about θ , the more we average over all possible parameter values.

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 - Probability Theory
 - Bayes Decision Theory
 - Maximum Likelihood Estimation
 - Bayesian Estimation
- **A Probabilistic View on Regression**
 - **Least-Squares Estimation as Maximum Likelihood**
 - **Predictive Distribution**
 - **Maximum-A-Posteriori (MAP) Estimation**
 - **Bayesian Curve Fitting**
- Discussion

Curve Fitting Revisited

- In the last lecture, we've looked at curve fitting in terms of error minimization...
- Now view the problem from a probabilistic perspective
 - Goal is to make predictions for target variable t given new value for input variable x .
 - Basis: training set $\mathbf{x} = (x_1, \dots, x_N)^T$ with target values $\mathbf{t} = (t_1, \dots, t_N)^T$.
 - We express our uncertainty over the value of the target variable using a probability distribution

$$p(t|x, \mathbf{w}, \beta)$$

Probabilistic Regression

- **First assumption:**

- Our target function values t are generated by adding noise to the ideal function estimate:

$$t = y(\mathbf{x}, \mathbf{w}) + \epsilon$$

Target function value

Regression function

Input value

Weights or parameters

Noise

- **Second assumption:**

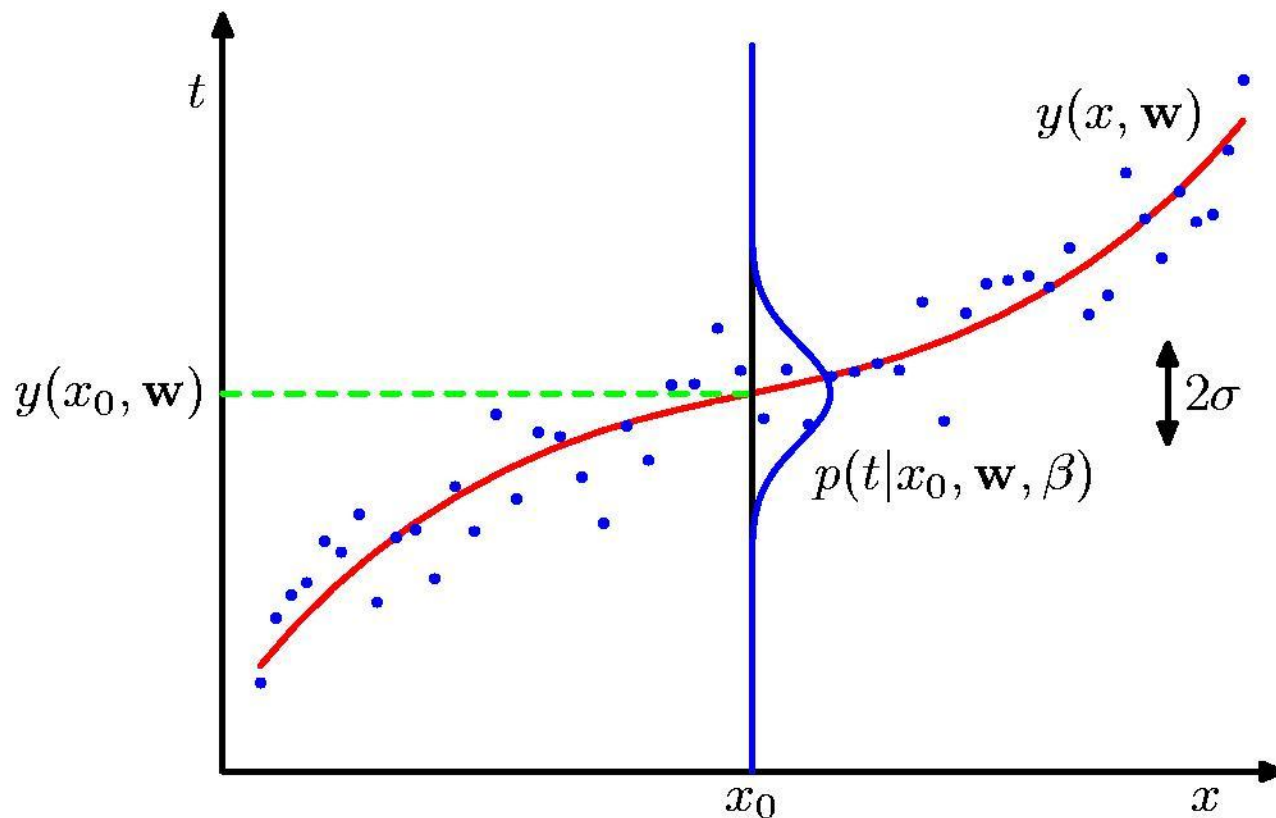
- The noise is Gaussian distributed.

$$p(t|\mathbf{x}, \mathbf{w}, \beta) = \mathcal{N}(t|y(\mathbf{x}, \mathbf{w}), \beta^{-1})$$

Mean

Variance
(β precision)

Visualization: Gaussian Noise



Probabilistic Regression

- **Given**

- Training data points:

$$\mathbf{X} = [\mathbf{x}_1, \dots, \mathbf{x}_n] \in \mathbb{R}^{d \times n}$$

- Associated function values:

$$\mathbf{t} = [t_1, \dots, t_n]^T$$

- **Conditional likelihood (assuming i.i.d. data)**

$$p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = \prod_{n=1}^N \mathcal{N}(t_n | y(\mathbf{x}_n, \mathbf{w}), \beta^{-1}) = \prod_{n=1}^N \mathcal{N}(t_n | \underbrace{\mathbf{w}^T \phi(\mathbf{x}_n)}_{\text{Generalized linear regression function}}, \beta^{-1})$$

⇒ Maximize w.r.t. \mathbf{w}, β

Generalized linear regression function

Maximum Likelihood Regression

- Simplify the log-likelihood

$$\log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = \sum_{n=1}^N \log \mathcal{N}(t_n | y(\mathbf{x}_n, \mathbf{w}), \beta^{-1})$$

$$\mathcal{N}(x|\mu, \beta^{-1}) = \frac{1}{\sqrt{2\pi}\beta^{-1/2}} \exp\left\{-\frac{\beta}{2}(x-\mu)^2\right\}$$

$$= \sum_{n=1}^N \left[\log\left(\frac{\sqrt{\beta}}{\sqrt{2\pi}}\right) - \frac{\beta}{2} \{y(\mathbf{x}_n, \mathbf{w}) - t_n\}^2 \right]$$

$$= -\frac{\beta}{2} \underbrace{\sum_{n=1}^N \{t_n - y(\mathbf{x}_n, \mathbf{w})\}^2}_{\text{Sum-of-squares error}} + \underbrace{\frac{N}{2} \log \beta - \frac{N}{2} \log(2\pi)}_{\text{Constants}}$$

Sum-of-squares error

Constants

Maximum Likelihood Regression

$$\begin{aligned}\log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) &= -\frac{\beta}{2} \sum_{n=1}^N \{t_n - y(\mathbf{x}_n, \mathbf{w})\}^2 + \frac{N}{2} \log \beta - \frac{N}{2} \log(2\pi) \\ &= -\frac{\beta}{2} \sum_{n=1}^N \{t_n - \mathbf{w}^T \phi(\mathbf{x}_n)\}^2 + \frac{N}{2} \log \beta - \frac{N}{2} \log(2\pi)\end{aligned}$$

- Gradient w.r.t. \mathbf{w} :

$$\nabla_{\mathbf{w}} \log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = -\beta \sum_{n=1}^N (t_n - \mathbf{w}^T \phi(\mathbf{x}_n)) \phi(\mathbf{x}_n)$$

Maximum Likelihood Regression

$$\nabla_{\mathbf{w}} \log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = -\beta \sum_{n=1}^N (t_n - \mathbf{w}^T \phi(\mathbf{x}_n)) \phi(\mathbf{x}_n)$$

- **Setting the gradient to zero:**

$$0 = -\beta \sum_{n=1}^N (t_n - \mathbf{w}^T \phi(\mathbf{x}_n)) \phi(\mathbf{x}_n)$$

$$\Leftrightarrow \sum_{n=1}^N y_i \phi(\mathbf{x}_n) = \left[\sum_{n=1}^N \phi(\mathbf{x}_n) \phi(\mathbf{x}_n)^T \right] \mathbf{w}$$

$$\Leftrightarrow \mathbf{\Phi} \mathbf{t} = \mathbf{\Phi} \mathbf{\Phi}^T \mathbf{w} \quad \mathbf{\Phi} = [\phi(\mathbf{x}_1), \dots, \phi(\mathbf{x}_n)]$$

$$\Leftrightarrow \mathbf{w}_{\text{ML}} = (\mathbf{\Phi} \mathbf{\Phi}^T)^{-1} \mathbf{\Phi} \mathbf{t} \quad \leftarrow \text{Same as in least-squares regression!}$$

\Rightarrow Least-squares regression is equivalent to Maximum Likelihood under the assumption of Gaussian noise.

Role of the Precision Parameter

- Also use ML to determine the precision parameter β :

$$\log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = -\frac{\beta}{2} \sum_{n=1}^N \{t_n - \mathbf{w}^T \phi(\mathbf{x}_n)\}^2 + \frac{N}{2} \log \beta - \frac{N}{2} \log(2\pi)$$

- Gradient w.r.t. β :

$$\nabla_{\beta} \log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = -\frac{1}{2} \sum_{n=1}^N \{t_n - \mathbf{w}^T \phi(\mathbf{x}_n)\}^2 + \frac{N}{2} \frac{1}{\beta}$$

$$\frac{1}{\beta_{\text{ML}}} = \frac{1}{N} \sum_{n=1}^N \{t_n - \mathbf{w}^T \phi(\mathbf{x}_n)\}^2$$

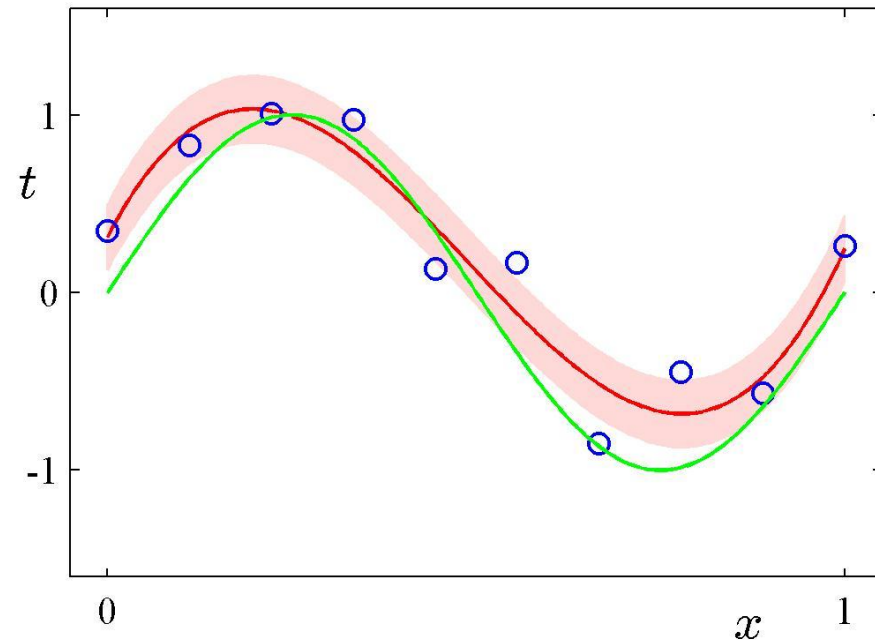
\Rightarrow *The inverse of the noise precision is given by the residual variance of the target values around the regression function.*

Predictive Distribution

- Having determined the parameters \mathbf{w} and β , we can now make predictions for new values of \mathbf{x} .

$$p(t|\mathbf{X}, \mathbf{w}_{\text{ML}}, \beta_{\text{ML}}) = \mathcal{N}(t|y(\mathbf{x}, \mathbf{w}_{\text{ML}}), \beta_{\text{ML}}^{-1})$$

- This means
 - Rather than giving a point estimate, we can now also give an estimate of the estimation uncertainty.



MAP: A Step Towards Bayesian Estimation...

- Introduce a prior distribution over the coefficients \mathbf{w} .

- For simplicity, assume a zero-mean Gaussian distribution

$$p(\mathbf{w}|\alpha) = \mathcal{N}(\mathbf{w}|\mathbf{0}, \alpha^{-1}\mathbf{I}) = \left(\frac{\alpha}{2\pi}\right)^{(M+1)/2} \exp\left\{-\frac{\alpha}{2}\mathbf{w}^T\mathbf{w}\right\}$$

- New **hyperparameter** α controls the distribution of model parameters.

- Express the posterior distribution over \mathbf{w} .

- Using Bayes' theorem:

$$p(\mathbf{w}|\mathbf{X}, \mathbf{t}, \beta, \alpha) \propto p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta)p(\mathbf{w}|\alpha)$$

- We can now determine \mathbf{w} by maximizing the posterior.
- This technique is called **maximum-a-posteriori (MAP)**.

MAP Solution

- Minimize the negative logarithm

$$-\log p(\mathbf{w}|\mathbf{X}, \mathbf{t}, \beta, \alpha) \propto -\log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) - \log p(\mathbf{w}|\alpha)$$

$$-\log p(\mathbf{t}|\mathbf{X}, \mathbf{w}, \beta) = \frac{\beta}{2} \sum_{n=1}^N \{y(\mathbf{x}_n, \mathbf{w}) - t_n\}^2 + \text{const}$$

$$-\log p(\mathbf{w}|\alpha) = \frac{\alpha}{2} \mathbf{w}^T \mathbf{w} + \text{const}$$

- The MAP solution is therefore

$$\frac{\beta}{2} \sum_{n=1}^N \{y(\mathbf{x}_n, \mathbf{w}) - t_n\}^2 + \frac{\alpha}{2} \mathbf{w}^T \mathbf{w}$$

\Rightarrow *Maximizing the posterior distribution is equivalent to minimizing the regularized sum-of-squares error (with $\lambda = \frac{\alpha}{\beta}$).*

Results of Probabilistic View on Regression

- Better understanding what linear regression *means*
 - *Least-squares regression is equivalent to ML estimation under the assumption of Gaussian noise.*
 - ⇒ We can use the **predictive distribution** to give an uncertainty estimate on the prediction.
 - ⇒ But: known problem with ML that it tends towards **overfitting**.
 - *L2-regularized regression (**Ridge regression**) is equivalent to MAP estimation with a Gaussian prior on the parameters w .*
 - ⇒ The prior controls the parameter values to reduce overfitting.
 - ⇒ This gives us a tool to explore more general priors.
- But still no full Bayesian Estimation yet
 - Should integrate over all values of w instead of point estimate.

Bayesian Curve Fitting

- **Given**

- Training data points: $\mathbf{X} = [\mathbf{x}_1, \dots, \mathbf{x}_n] \in \mathbb{R}^{d \times n}$
- Associated function values: $\mathbf{t} = [t_1, \dots, t_n]^T$
- Our goal is to predict the value of t for a new point \mathbf{x} .

- **Evaluate the predictive distribution**

$$p(t|x, \mathbf{X}, \mathbf{t}) = \int \underbrace{p(t|x, \mathbf{w})}_{\text{Noise distribution}} \underbrace{p(\mathbf{w}|\mathbf{X}, \mathbf{t})}_{\text{What we just computed for MAP}} d\mathbf{w}$$

What we just computed for MAP

- Noise distribution - again assume a Gaussian here

$$p(t|x, \mathbf{w}) = \mathcal{N}(t|y(\mathbf{x}, \mathbf{w}), \beta^{-1})$$

- Assume that parameters α and β are fixed and known for now.

Bayesian Curve Fitting

- Under those assumptions, the posterior distribution is a Gaussian and can be evaluated analytically:

$$p(t|x, \mathbf{X}, \mathbf{t}) = \mathcal{N}(t|m(x), s^2(x))$$

- where the mean and variance are given by

$$m(x) = \beta \phi(x)^T \mathbf{S} \sum_{n=1}^N \phi(\mathbf{x}_n) t_n$$

$$s(x)^2 = \beta^{-1} + \phi(x)^T \mathbf{S} \phi(x)$$

- and \mathbf{S} is the regularized covariance matrix

$$\mathbf{S}^{-1} = \alpha \mathbf{I} + \beta \sum_{n=1}^N \phi(\mathbf{x}_n) \phi(\mathbf{x}_n)^T$$

Analyzing the result

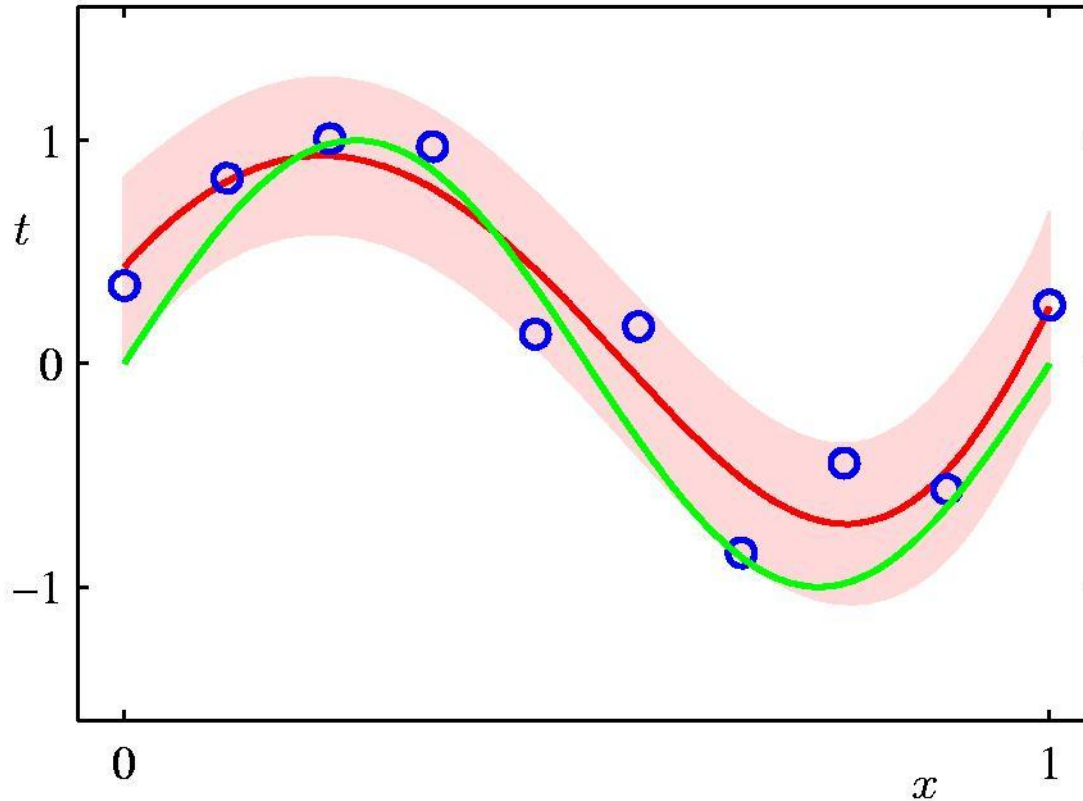
- Analyzing the variance of the predictive distribution

$$s(x)^2 = \underbrace{\beta^{-1}}_{\text{Uncertainty in the predicted value due to noise on the target variables (expressed already in ML)}} + \underbrace{\phi(x)^T \mathbf{S} \phi(x)}_{\text{Uncertainty in the parameters } w \text{ (consequence of Bayesian treatment)}}$$

Uncertainty in the predicted value due to noise on the target variables (expressed already in ML)

Uncertainty in the parameters w (consequence of Bayesian treatment)

Bayesian Predictive Distribution



- Important difference to previous example
 - Uncertainty may vary with test point x !

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 - Bayesian Curve Fitting
- **Discussion**

Discussion

- We now have a better understanding of regression
 - Least-squares regression: Assumption of Gaussian noise
 - ⇒ We can now also plug in different noise models and explore how they affect the error function.
 - L2 regularization as a Gaussian prior on parameters w .
 - ⇒ We can now also use different regularizers and explore what they mean.
 - ⇒ Next lecture...
 - General formulation with basis functions $\phi(\mathbf{x})$.
 - ⇒ We can now also use different basis functions.

Discussion

- **General regression formulation**
 - In principle, we can perform regression in arbitrary spaces and with many different types of basis functions
 - However, there is a caveat... Can you see what it is?

- **Example: Polynomial curve fitting, $M = 3$**

$$y(\mathbf{x}, \mathbf{w}) = w_0 + \sum_{i=1}^D w_i x_i + \sum_{i=1}^D \sum_{j=1}^D w_{ij} x_i x_j + \sum_{i=1}^D \sum_{j=1}^D \sum_{k=1}^D w_{ijk} x_i x_j x_k$$

⇒ Number of coefficients grows with D^M !

⇒ The approach becomes quickly unpractical for high dimensions.

- This is known as the **curse of dimensionality**.
- We will encounter some ways to deal with this later.

References and Further Reading

- More information on linear regression can be found in Chapters 1.2.5-1.2.6 and 3.1-3.1.4 of

Christopher M. Bishop
Pattern Recognition and Machine Learning
Springer, 2006

