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# Advanced Machine Learning Lecture 2

## Linear Regression

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## This Lecture: Advanced Machine Learning

- Regression Approaches
  - Linear Regression
  - Regularization (Ridge, Lasso)
  - Gaussian Processes
- Learning with Latent Variables
  - EM and Generalizations
  - Approximate Inference
- Deep Learning
  - Neural Networks
  - CNNs, RNNs, RBMs, etc.

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## 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

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

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## 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:
 

a a b a b a a b a  
 b a a a b a a b a  
 a b a a a b b a  
 b a b a a b a

?

$C_1 = a$   $p(C_1) = 0.75$

$C_2 = b$   $p(C_2) = 0.25$
- In general:  $\sum_k p(C_k) = 1$

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## 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  $C_k$ .

$p(x|a)$

$p(x|b)$

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## 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**

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

Slide credit: Bernt Schiele B. Leibe 7

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## Recap: Bayes Decision Theory

Likelihood

Likelihood  $\times$  Prior

Decision boundary

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

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## Recap: Gaussian (or Normal) Distribution

- **One-dimensional case**
  - Mean  $\mu$
  - Variance  $\sigma^2$
$$\mathcal{N}(x|\mu, \sigma^2) = \frac{1}{\sqrt{2\pi}\sigma} \exp\left\{-\frac{(x-\mu)^2}{2\sigma^2}\right\}$$
- **Multi-dimensional case**
  - Mean  $\mu$
  - Covariance  $\Sigma$
$$\mathcal{N}(x|\mu, \Sigma) = \frac{1}{(2\pi)^{D/2} |\Sigma|^{1/2}} \exp\left\{-\frac{1}{2}(x-\mu)^T \Sigma^{-1} (x-\mu)\right\}$$

Image source: C. M. Bishop, 2006 B. Leibe 9

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## Side Note

- **Notation**
  - In many situations, it will be preferable to work with the inverse of the covariance matrix  $\Sigma$ :
$$\Lambda = \Sigma^{-1}$$
  - We call  $\Lambda$  the **precision matrix**.
  - We can therefore also write the Gaussian as
$$\mathcal{N}(x|\mu, \Lambda^{-1}) = \frac{1}{\sqrt{2\pi} \lambda^{-1/2}} \exp\left\{-\frac{\lambda}{2}(x-\mu)^2\right\}$$

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

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## Recap: Parametric Methods

- **Given**
  - Data  $X = \{x_1, x_2, \dots, x_N\}$
  - Parametric form of the distribution with parameters  $\theta$
  - E.g. for Gaussian distrib.:  $\theta = (\mu, \sigma)$
- **Learning**
  - Estimation of the parameters  $\theta$
- **Likelihood of  $\theta$** 
  - Probability that the data  $X$  have indeed been generated from a probability density with parameters  $\theta$ 

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

Slide adapted from Bernt Schiele B. Leibe 11

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## Recap: Maximum Likelihood Approach

- **Computation of the likelihood**
  - Single data point:  $p(x_n|\theta) = \mathcal{N}(x_n|\mu, \sigma^2)$
  - 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  $\theta$  (Learning)**
  - Maximize the likelihood (=minimize the negative log-likelihood)
    - ⇒ Take the derivative and set it to zero.
$$\frac{\partial}{\partial \theta} E(\theta) = -\sum_{n=1}^N \frac{\partial}{\partial \theta} \ln p(x_n|\theta) \stackrel{!}{=} 0$$

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## 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...

13

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## 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.

14

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## 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...

15

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## Recap: Bayesian Approach to Learning

- Conceptual shift
  - Maximum Likelihood views the true parameter vector  $\theta$  to be unknown, but fixed.
  - In Bayesian learning, we consider  $\theta$  to be a random variable.
- This allows us to use knowledge about the parameters  $\theta$ 
  - i.e. to use a prior for  $\theta$
  - Training data then converts this prior distribution on  $\theta$  into a posterior probability density.

- The prior thus encodes knowledge we have about the type of distribution we expect to see for  $\theta$ .

16

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## Recap: Bayesian Learning Approach

- Bayesian view:**
  - Consider the parameter vector  $\theta$  as a random variable.
  - When estimating the parameters, what we compute is

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

Assumption: given  $\theta$ , this doesn't depend on  $X$  anymore

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

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

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

17

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## 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

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

18

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## Recap: Bayesian Learning Approach

- Discussion
  - Likelihood of the parametric form  $\theta$  given the data set  $X$ .
  - Prior for the parameters  $\theta$
  - Estimate for  $x$  based on parametric form  $\theta$

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

Normalization: integrate over all possible values of  $\theta$

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

19

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  - 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

20

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## 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

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

21

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## 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  $\leftarrow$   $t = y(\mathbf{x}, \mathbf{w}) + \epsilon$   $\leftarrow$  Noise  
 Regression function  $\leftarrow$   $y(\mathbf{x}, \mathbf{w})$   $\leftarrow$  Input value  $\leftarrow$  Weights or parameters

- Second assumption:
  - The noise is Gaussian distributed.

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

Mean  $\leftarrow$   $y(\mathbf{x}, \mathbf{w})$   $\leftarrow$  Variance ( $\beta$  precision)

22

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## Visualization: Gaussian Noise

23

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

Generalized linear regression function

$\Rightarrow$  Maximize w.r.t.  $\mathbf{w}, \beta$

24

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## Maximum Likelihood Regression

- Simplify the log-likelihood

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

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

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

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

Slide adapted from Bernt Schiele. B. Leibe. 25

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## Maximum Likelihood Regression

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

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

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

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

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## 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:

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

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

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

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

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

Slide adapted from Bernt Schiele. B. Leibe. 28

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## Role of the Precision Parameter

- Also use ML to determine the precision parameter  $\beta$ :

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

- Gradient w.r.t.  $\beta$ :

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

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

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

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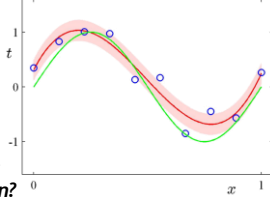
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## Predictive Distribution

- Having determined the parameters  $\mathbf{w}$  and  $\beta$ , 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.



- What else can we do in the Bayesian view of regression?

B. Leibe. Image source: C.M. Bishop, 2006. 30

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## 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  $\alpha$  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)**.

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## 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 the solution of
  - $\frac{\beta}{2} \sum_{n=1}^N \{y(\mathbf{x}_n, \mathbf{w}) - t_n\}^2 + \frac{\alpha}{2} \mathbf{w}^T \mathbf{w}$
  - $\Rightarrow$  Maximizing the posterior distribution is equivalent to minimizing the regularized sum-of-squares error (with  $\lambda = \frac{\alpha}{\beta}$ ).

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## 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.
  - $\Rightarrow$  We can use the predictive distribution to give an uncertainty estimate on the prediction.
  - $\Rightarrow$  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  $\mathbf{w}$ .
  - $\Rightarrow$  The prior controls the parameter values to reduce overfitting.
  - $\Rightarrow$  This gives us a tool to explore more general priors.
- But still no full Bayesian Estimation yet
  - Should integrate over all values of  $\mathbf{w}$  instead of just making a point estimate.

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## 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
  - $$p(t|\mathbf{x}, \mathbf{X}, \mathbf{t}) = \int p(t|\mathbf{x}, \mathbf{w}) p(\mathbf{w}|\mathbf{X}, \mathbf{t}) d\mathbf{w}$$
  - What we just computed for MAP
  - Noise distribution - again assume a Gaussian here
    - $p(t|\mathbf{x}, \mathbf{w}) = \mathcal{N}(t|y(\mathbf{x}, \mathbf{w}), \beta^{-1})$
  - Assume that parameters  $\alpha$  and  $\beta$  are fixed and known for now.

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## Bayesian Curve Fitting

- Under those assumptions, the posterior distribution is a Gaussian and can be evaluated analytically:
  - $$p(t|\mathbf{x}, \mathbf{X}, \mathbf{t}) = \mathcal{N}(t|m(x), s^2(x))$$
  - 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
    - $$\mathbf{S}^{-1} = \alpha \mathbf{I} + \beta \sum_{n=1}^N \phi(\mathbf{x}_n) \phi(\mathbf{x}_n)^T$$

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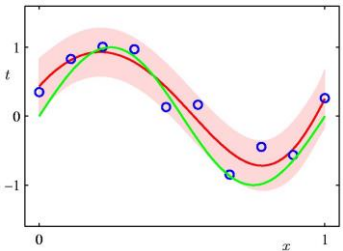
## Analyzing the result

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

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## Bayesian Predictive Distribution



- Important difference to previous example
  - Uncertainty may vary with test point  $\mathbf{x}$ !

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38

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## 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  $\phi(x)$ .
    - ⇒ We can now also use different basis functions.

39

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## 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.

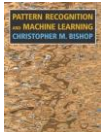
40

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## 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



41