# Advanced probabilistic methods Lecture 3: Multivariate Gaussian, Bayesian linear models

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- Gaussian distribution
  - Bayesian parameter learning
- Multivariate Gaussian distribution
  - Characterization
  - Useful identities
- Bayesian Linear Parameter Models (LPMs)
  - Posterior computation (given fixed hyperparameters)
- Ch. 8 & 18 (until the end of Section 18.1.1) in Barber's book

# Recall from lecture 1

- Tools for probabilistic modeling
  - **Models**: Bayesian networks, sparse Bayesian linear regression, Gaussian mixture models, latent linear models
  - Methods for inference: maximum likelihood, maximum a posteriori (MAP), analytical, Laplace approximation, expectation maximization (EM), Variational Bayes (VB), stochastic variational inference (SVI)
  - Ways to select between models



- A model specifies a probability distribution for a random variable Y, and it is often affected by some parameter θ. The model can be denoted as p(y|θ).
- Fitting the model (i.e. inference) corresponds to learning the value (or the distribution) of θ, after some data y have been observed.

# Prior, Likelihood, Posterior

 Bayes' rule tells us how to update our prior beliefs about variable θ in light of the data y to a posterior belief:



The evidence is also called the marginal likelihood.

- $p(y|\theta)$  is the probability that the model generates the observed data y when using parameter  $\theta$ 
  - $L(\theta) \equiv p(y|\theta)$ , with y held fixed, is called the *likelihood*
  - $f(y) \equiv p(y|\theta)$ , with  $\theta$  held fixed, is called the *observation model*
- "*Methods for inference*" = Bayes' rule + some algorithm to do the actual computations (on this course)

• The *Maximum A Posteriori (MAP)* parameter value, which maximizes the posterior

$$heta_* = rg\max_{ heta} p( heta|y)$$

• The Maximum likelihood assignment (ML)

$$heta_* = rg\max_{ heta} p(y| heta)$$

• The full posterior distribution  $p(\theta|y)$  tells also of the uncertainty about the value of  $\theta$ .

# Gaussian distribution

- $X \sim N(\mu, \sigma^2)$
- Parameters:  $\mu$ : mean,  $\sigma^2$ : variance
- Inverse of the variance,  $\lambda=1/\sigma^2$ , is called the precision
- Standard deviation  $\sigma$
- 95% credible interval equals approximately  $[\mu-2\sigma,\mu+2\sigma]$

PDF:

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



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# Bayesian estimation of the mean of a Gaussian (1/2)

- Suppose we have observations  $x = (x_1, ..., x_n)$  from  $N(\mu, \sigma^2)$ , where  $\sigma^2$  is known.
- To learn  $\mu$ , we specify a prior

$$\mu \sim N(\mu_0, au_0^2)$$

Posterior

$$p(\mu|x) = \frac{p(x|\mu)p(\mu)}{p(x)} \propto p(\mu)p(x|\mu)$$
  
=  $\frac{1}{\sqrt{2\pi\tau_0}} e^{-\frac{1}{2\tau_0^2}(\mu-\mu_0)^2} \times \prod_{i=1}^n \frac{1}{\sqrt{2\pi\sigma}} e^{-\frac{1}{2\sigma^2}(x_i-\mu)^2}$   
 $\propto e^{-\frac{1}{2\tau_0^2}(\mu-\mu_0)-\frac{1}{2\sigma^2}\sum_i(x_i-\mu)^2}$   
= ... (details in BDA course)

Bayesian estimation of the mean of a Gaussian (2/2)

Posterior

$$p(\mu|x) \propto e^{-\frac{1}{2\tau_n^2}(\mu-\mu_n)^2}$$
$$\propto N(\mu|\mu_n, \tau_n^2)$$

where

$$\mu_n = \frac{\frac{1}{\tau_0^2}\mu_0 + \frac{n}{\sigma^2}\overline{x}}{\frac{n}{\sigma^2} + \frac{1}{\tau_0^2}} \quad \text{and} \quad \frac{1}{\tau_n^2} = \frac{n}{\sigma^2} + \frac{1}{\tau_0^2}$$

• Posterior precision  $1/\tau_n^2$ : sum of prior precision  $1/\tau_0^2$  and data precision  $n/\sigma^2$ 

 Posterior mean μ<sub>n</sub>: precision weighted average of prior mean μ<sub>0</sub> and data mean x̄. • In the previous example

Prior: 
$$\mu \sim N(\mu_0, \tau_0^2)$$
  
Posterior:  $\mu \sim N(\mu_n, \tau_n^2)$ .

If the prior and posterior belong to the same family of distributions, we say that the prior is conjugate to the likelihood used.

• For example, normal prior  $\mu \sim N(\mu_0, \tau_0^2)$  is conjugate to the normal likelihood  $N(x|\mu, \sigma^2)$ .

• Conjugacy is useful, because it makes computations easy.

• With conjugate prior, the posterior is available in a closed form

 $p(\boldsymbol{\theta}|\boldsymbol{x}) \propto p(\boldsymbol{x}|\boldsymbol{\theta})p(\boldsymbol{\theta})$ 

- Drop all terms not depending on  $\boldsymbol{\theta}$
- Recognize the result as a density function belonging to the same family of distributions as the prior  $p(\theta)$ , but with different parameters.
- Examples (likelihood conjugate prior):
  - Likelihood for normal mean Normal prior
  - Likelihood for normal variance Inverse-Gamma prior
  - Bernoulli Beta
  - Binomial Beta
  - Exponential Gamma
  - Poisson Gamma

- Suppose we have observations  $x = (x_1, ..., x_n)$  from  $N(\mu, \lambda^{-1})$ , where both the mean  $\mu$  and the precision  $\lambda$  are unknown.
- The conjugate prior distribution is the normal-gamma distribution

$$p(\mu, \lambda | \mu_0, \beta, a, b) = N(\mu | \mu_0, (\beta \lambda)^{-1}) \operatorname{Gam}(\lambda | a, b)$$
  
 $\equiv \operatorname{Normal-Gamma}(\mu, \lambda | \mu_0, \beta, a, b)$ 

Note the dependency of the prior of  $\mu$  on the value of  $\lambda$ .

# Gaussian distribution, unknown mean and precision (2/2)

• The conjugate prior distribution is the normal-gamma distribution  $p(\mu, \lambda | \mu_0, \beta, a, b) =$ Normal-Gamma $(\mu, \lambda | \mu_0, \beta, a, b)$ 

Posterior

$$p(\mu, \lambda | x) = \mathsf{Normal-Gamma}(\mu, \lambda | \mu_n, \beta_n, a_n, b_n),$$

with

$$\mu_n = \frac{\beta \mu_0 + n\overline{x}}{\beta + n}$$
  

$$\beta_n = \beta + n$$
  

$$a_n = a + \frac{n}{2}$$
  

$$b_n = b + \frac{1}{2} \left( ns + \frac{\beta n(\overline{x} - \mu_0)^2}{\beta + n} \right)$$

# Gaussian distribution, unknown mean and precision, example (1/2)

- Simulate samples from  $N(\mu = 2, \sigma^2 = 0.25)$ 
  - precision  $\lambda = 4$
- Try to learn  $\mu$  and  $\lambda$
- Specify prior

$$m{
ho}(\mu,\lambda|\mu_0,eta,m{a},m{b})=\mathsf{Normal-Gamma}(\mu,\lambda|\mu_0,m{eta},m{a},m{b})$$

with

$$\mu_0=$$
 0,  $\beta=0.001,$   $a=0.01,$   $b=0.01$ 

• See: normal\_example.m

# Gaussian distribution, unknown mean and precision, example (2/2)

• When  $\mu$  and  $\lambda$  have distribution

Normal-Gamma $(\mu, \lambda | \mu_n, \beta_n, a_n, b_n) = N(\mu | \mu_n, (\beta_n \lambda)^{-1})$ Gam $(\lambda | a_n, b_n)$ ,

marginal distribution of  $\lambda$  can be plotted using the PDF of  ${\rm Gam}(\lambda|\mathbf{a}_n,\mathbf{b}_n)$ 

- To plot the marginal distribution of  $\mu$ , we need to take the dependence on  $\lambda$  into account.
  - we compute the marginal distribution of  $\mu$  by averaging over  $N(\mu|\mu_n, (\beta_n\lambda_i)^{-1})$ , for multiple  $\lambda_i$  simulated from  $Gam(\lambda|a_n, b_n)$
  - (could also be done analytically...)

• If  $p(x|\theta_t)$  is the true data generating mechanism, and A is a neighborhood of  $\theta_t$ , then

$$p(\theta \in A|x) \stackrel{n \to \infty}{\to} 1.$$

- The posterior distribution concentrates around the true value (if such a value exists!). See the *normal\_example.m*
- It follows that

$$\overline{\theta}_{MAP} \stackrel{n \to \infty}{\to} \theta_t$$
 and  $\overline{\theta}_{ML} \stackrel{n \to \infty}{\to} \theta_t$ 

$$N_D(x|\mu, \Sigma) \equiv (2\pi)^{-\frac{D}{2}} |\Sigma|^{-\frac{1}{2}} e^{-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)}$$

• D: dimension,  $\mu$ : mean,  $\Sigma$ : covariance matrix. With D = 2:

$$\mu = \begin{bmatrix} \mu_1 \\ \mu_2 \end{bmatrix}, \quad \Sigma = \begin{bmatrix} \sigma_1^2 & \sigma_{12} \\ \sigma_{21} & \sigma_2^2 \end{bmatrix}$$

- σ<sub>12</sub> = σ<sub>21</sub>: covariance between x<sub>1</sub> and x<sub>2</sub>. (tells direction of dependency)
- $\rho_{12} = \sigma_{12}/(\sigma_1\sigma_2)$ :correlation between  $x_1$  and  $x_2$ . (direction and strength)



# Multivariate Gaussian - characterization (1/2)



Eigendecomposition

$$\Sigma = E \Lambda E^T$$
,

where  $E^T E = I$  and  $\Lambda = \text{diag}(\lambda_1, \ldots, \lambda_D)$ .

Now the transformation

$$y = \Lambda^{-\frac{1}{2}} E^T (x - \mu)$$

can be shown to have the distribution  $N_D(0, I)$  (product of D independent standard Gaussians)

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# Multivariate Gaussian - characterization (2/2)



- Thus,  $x = E\Lambda^{\frac{1}{2}}y + \mu$  with distribution  $N_D(\mu, \Sigma)$  is obtained from standard independent Gaussians y by
  - scaling by the square roots of eigenvalues
  - rotating by the eigenvectors
  - shifting by adding the mean

#### • Let $z \sim N(\mu, \Sigma)$ and consider partitioning it as:

$$z = \left(\begin{array}{c} x \\ y \end{array}\right)$$

with

$$\mu = \left( egin{array}{c} \mu_x \ \mu_y \end{array} 
ight)$$
 and  $\Sigma = \left( egin{array}{cc} \Sigma_{xx} & \Sigma_{xy} \ \Sigma_{yx} & \Sigma_{yy} \end{array} 
ight).$ 

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# Marginalization and conditioning (2/2)

• Then

$$\begin{split} p(x) &\sim \mathcal{N}(\mu_x, \Sigma_{xx}) \quad \text{(marginalization)} \\ p(x|y) &= \mathcal{N}(\mu_x + \Sigma_{xy} \Sigma_{yy}^{-1} (y - \mu_y), \Sigma_{xx} - \Sigma_{xy} \Sigma_{yy}^{-1} \Sigma_{yx}) \quad \text{(conditioning} \\ \Longrightarrow \text{Marginals and conditionals of M-V Gaussians are still M-V} \end{split}$$

Gaussian.



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• Linear transformation: if

$$y = Mx + \eta$$
,

where  $x \sim \textit{N}(\mu_x, \Sigma_x)$  and  $\eta \sim \textit{N}(\mu, \Sigma)$ ,then

$$p(y) = N(y|M\mu_x + \mu, M\Sigma_x M^T + \Sigma)$$

• Completing the square:

$$\frac{1}{2}x^{T}Ax - b^{T}x = \frac{1}{2}(x - A^{-1}b)^{T}A(x - A^{-1}b) - \frac{1}{2}b^{T}A^{-1}b$$

From which one can derive, for example

$$\int \exp(-\frac{1}{2}x^{\mathsf{T}}Ax + b^{\mathsf{T}}x)dx = \sqrt{\det(2\pi A^{-1})}\exp(\frac{1}{2}b^{\mathsf{T}}A^{-1}b)$$

Let x = (x<sub>1</sub>,..., x<sub>n</sub>) be from N(μ, Σ) with unknown μ and Σ.
 Log-likelihood, assuming data are *i.i.d.*:

$$\begin{split} L(\mu,\Sigma) &= \sum_{i=1}^{N} \log p(x_i | \mu, \Sigma) \\ &= -\frac{1}{2} \sum_{i=1}^{N} (x_i - \mu)^T \Sigma^{-1} (x_i - \mu) - \frac{N}{2} \log \det(2\pi\Sigma) \end{split}$$

#### Multivariate Gaussian - ML fitting

• Differentiate  $L(\mu, \Sigma)$  w.r.t. the vector  $\mu$ :

$$\nabla_{\mu} L(\mu, \Sigma) = \sum_{i=1}^{N} \Sigma^{-1}(x_i - \mu)$$

Equating to zero gives

$$\sum_{i=1}^N \Sigma^{-1} x_i = N \Sigma^{-1} \mu.$$

Thus we get

$$\widehat{\mu} = \frac{1}{N} \sum_{i=1}^{N} x_i$$

• Similarly one can derive:

$$\widehat{\Sigma} = \frac{1}{N} \sum_{i=1}^{N} (x_i - \overline{x}) (x_i - \overline{x})^T$$

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• Gaussian-Wishart is the conjugate prior, when  $X_i \sim N(\mu, \Lambda)$  and both mean  $\mu$  and precision  $\Lambda$  are unknown:

$$p(\mu, \Lambda | \mu_0, \beta, W, \nu) = N(\mu | \mu_0, (\beta \Lambda)^{-1}) \mathcal{W}(\Lambda | W, \nu)$$

If X<sub>i</sub> are scalar, this is equivalent to the Gaussian-Gamma distribution.
Posterior

$$p(\mu, \Lambda | x) = N(\mu | \mu_n, (\beta_n \Lambda)^{-1}) \mathcal{W}(\Lambda | W_n, \nu_n)$$

• Wishart distribution is a distribution for nonnegative-definite matrix-valued random variables

 $\Lambda \sim \mathcal{W}(\Lambda | \mathcal{W}, \nu)$ 

$$E(\Lambda) = 
u W$$
  
 $\mathsf{Var}(\Lambda_{ij}) = n(w_{ij}^2 + w_{ii}w_{jj})$ 

• Further: exercises...

#### Linear models with Gaussian noise

• Data 
$$\mathcal{D} = \{(x_i, y_i), i = 1, ..., N\}$$

- **x**<sub>i</sub>: the input
- *y<sub>i</sub>*: the output
- Model:

$$y = \underbrace{f(\mathbf{w}, \mathbf{x})}_{\text{clean output}} + \underbrace{\eta}_{\text{noise}}, \quad \eta \sim N(0, \beta^{-1})$$

• In the simplest case

$$f(\mathbf{w}, \mathbf{x}) = \mathbf{w}^T \mathbf{x}$$
$$= w_1 x_1 + \ldots + w_D x_D$$

• The parameters w<sub>i</sub> are also called the weights

- A prior distribution  $p(\mathbf{w}|\alpha)$  is placed on the weights  $\mathbf{w}$ .
- The posterior distribution  $p(\mathbf{w}|\mathcal{D}, \Gamma)$  can be computed, and reflects the uncertainty of the parameters.

• A Gaussian prior distribution may placed on w:

$$p(\mathbf{w}|\alpha) = N(\mathbf{w}|\mathbf{0}, \alpha^{-1}\mathbf{I})$$
$$= \prod_{i=1}^{D} N(w_i|\mathbf{0}, \alpha^{-1}) = \left(\frac{\alpha}{2\pi}\right)^{\frac{D}{2}} e^{-\frac{\alpha}{2}\sum_i w_i^2}$$

Posterior

$$\log p(\mathbf{w}|\Gamma, \mathcal{D}) = -\frac{\beta}{2} \sum_{i=1}^{N} \left[ y_i - \mathbf{w}^T \mathbf{x}_i \right]^2 - \frac{\alpha}{2} \mathbf{w}^T \mathbf{w} + \text{const}$$

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- *α*: precision of the regression weights
  - determines the amount of regularization
  - $\bullet~$  large precision  $\rightarrow$  small variance  $\rightarrow$  weights are close to zero
- $\beta$ : *precision* of the noise
- $\Gamma = \{\alpha, \beta\}$  are called the **hyperparameters** (in the course book...)

# Posterior distribution

Posterior distribution is obtained by completing the square (left as an exercise):

$$p(\mathbf{w}|\Gamma, \mathcal{D}) = N(\mathbf{w}|\mathbf{m}, S)$$

where

$$S = \left(\alpha I + \beta \sum_{i=1}^{N} \mathbf{x}_{i} \mathbf{x}_{i}^{T}\right)^{-1}, \qquad \mathbf{m} = \beta S \sum_{i=1}^{N} y_{i} \mathbf{x}_{i}$$

Mean prediction

$$\widetilde{y} = \int \mathbf{w}^T \mathbf{x} imes p(\mathbf{w} | \Gamma, \mathcal{D}) d\mathbf{w} = \mathbf{m}^T \mathbf{x}$$

- Setup: simulate  $y = \mathbf{w}_{true}^T \mathbf{x} + \epsilon$ , where  $\epsilon \sim N(0, \beta^{-1})$  and  $\beta = 1$
- The goal is to investigate how hyperparameter *α* affects the posterior distribution of the parameters w

#### Example, impact of hyperparameters (2/3)

• Too large  $\alpha$ ,  $Var(y - \widetilde{y}) = 1.54$  (Original Var(y) = 1.75)



• Too small  $\alpha$ ,  $Var(y - \tilde{y}) = 2.48$ 



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Advanced probabilistic methods

## Example, impact of hyperparameters (3/3)

- About good  $\alpha$ ,  $Var(y \tilde{y}) = 1.46$
- A compromise between bias and variance



• Other sparse priors (e.g., Laplace, horse-shoe, spike-and-slab):



- Analysis of  $\sim$  1,000,000 genetic polymorphisms in  $\sim$  50,000 genomic regions (Peltola et al., 2012, *PLoS ONE*).
- Spike-and-slab prior on regression weights



- Bayesian learning of the Gaussian distribution using conjugate priors
- Multivariate Gaussian
  - Characterization
  - Marginal & conditional distributions
  - Linear transformation & completing the square
- By placing a Gaussian prior on the parameters of linear regression, the posterior is also Gaussian.
- Meaning and impact of hyperparameters in Bayesian linear regression.