

# Deep Learning - Foundations and Concepts

## Chapter 2. Probabilities

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# Outline

- 1 The Rules of Probability
- 2 Probability Densities
- 3 The Gaussian Distribution
- 4 Transformation of Densities
- 5 Information Theory
- 6 Bayesian Probabilities

# The sum and product rules

- Sum rule:  $p(X) = \sum_Y p(X, Y)$ .
- Product rule:  $p(X, Y) = p(Y|X)p(X)$ .

# Bayes' theorem

- Bayes' theorem:

$$\begin{aligned} p(Y|X) &= \frac{p(X|Y)p(Y)}{p(X)} \\ &= \frac{p(X|Y)p(Y)}{\sum_Y p(X|Y)p(Y)} \end{aligned}$$

- Prior and posterior probabilities:
  - $p(Y)$  is the prior probability, because it is available *before* we observe the event  $X$ .
  - $p(Y|X)$  is the posterior probability, because it is obtained *after* we have observed the event  $X$ .

# Probability densities

- A probability density  $p(x)$  is a real function satisfies the following two conditions<sup>1</sup>:
  - $p(x) \geq 0$ .
  - $\int_{-\infty}^{+\infty} p(x)dx = 1$ .
- The *cumulative distribution function* is given by  $P(x) = \int_{-\infty}^x p(t)dt$ , and usually we have  $P'(x) = p(x)$ .
- These definitions can easily be extended to higher dimensions.

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<sup>1</sup>This is *not* a mathematically robust definition, but it suffices for this course.

# Probability densities

- Sum rule:  $p(x) = \int p(x, y)dy$ .
- Product rule:  $p(x, y) = p(y|x)p(x)$ .
- Bayes' theorem:  $p(y|x) = \frac{p(x|y)p(y)}{p(x)} = \frac{p(x|y)p(y)}{\int p(x|y)p(y)dy}$ .

# Expectations and covariances

- Expectation of  $f$ :
  - Discrete case:  $E(f) = \sum_x p(x)f(x)$ .
  - Continuous case:  $E(f) = \int p(x)f(x)dx$ .
- Variance of  $f$ :  $\text{var}(f) = E((f(x) - E(f))^2) = E(f^2) - E(f)^2$ .
- Covariance of:
  - Two random variables:
 
$$\text{cov}(x, y) = E((x - E(x))(y - E(y))) = E(xy) - E(x)E(y).$$
  - Two vectors:
 
$$\text{cov}(x, y) = E((x - E(x))(y - E(y))^T) = E(xy^T) - E(x)E(y^T).$$

# Example distributions

- Uniform distribution:  $p(x) = \frac{1}{d-c}, \quad x \in (c, d).$
- Exponential distribution:  $p(x; \lambda) = \lambda \exp(-\lambda x).$
- Laplace distribution:  $p(x; \mu, \gamma) = \frac{1}{2\gamma} \exp(-\frac{|x-\mu|}{\gamma}).$
- Dirac delta function:  $p(x; \mu_1, \dots, \mu_N) = \frac{1}{N} \sum_{n=1}^N \delta(x - \mu_n).$



# The Gaussian distribution

- Definition:  $\mathcal{N}(x; \mu, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp(-\frac{(x-\mu)^2}{2\sigma^2})$ .
- Mean:  $E(x) = \int_{-\infty}^{+\infty} \mathcal{N}(x; \mu, \sigma^2) x dx = \mu$ .
- Variance:  $\text{var}(x) = E(x^2) - E(x)^2 = \sigma^2$ .

# Maximum likelihood and its bias

## Problem

We have  $N$  observations of a random variable  $x$ :  $x_1, \dots, x_N$  that are drawn independently from a Gaussian distribution whose mean  $\mu$  and variance  $\sigma^2$  are unknown. How do we determine these parameters from the data set?

# Maximum likelihood and its bias

## Problem'

Find  $\mu$  and  $\sigma^2$  such that the probability of the data set

$$p(x_1, \dots, x_N; \mu, \sigma^2) = \prod_{n=1}^N \mathcal{N}(x_n; \mu, \sigma^2)$$

is maximized.

# Maximum likelihood and its bias

Problem"

Let's minimize

$$\begin{aligned} L &= -\log p(x_1, \dots, x_N; \mu, \sigma^2) \\ &= \frac{1}{2\sigma^2} \sum_{n=1}^N (x_n - \mu)^2 + \frac{N}{2} \log \sigma^2 + \frac{N}{2} \log(2\pi) \end{aligned}$$

instead.

# Maximum likelihood and its bias

$$\frac{\partial L}{\partial \mu} = \frac{1}{\sigma^2} \sum_{n=1}^N (\mu - x_n) = \frac{N}{\sigma^2} \left( \mu - \frac{1}{N} \sum_{n=1}^N x_n \right)$$

$$\frac{\partial L}{\partial \sigma} = \frac{N}{\sigma^3} \left( \sigma^2 - \frac{1}{N} \sum_{n=1}^N (x_n - \mu)^2 \right)$$

Setting  $\frac{\partial L}{\partial \mu}$  and  $\frac{\partial L}{\partial \sigma}$  to 0, we have:

$$\mu_{ML} = \frac{1}{N} \sum_{n=1}^N x_n$$

$$\sigma_{ML}^2 = \frac{1}{N} \sum_{n=1}^N (x_n - \mu_{ML})^2$$

# Maximum likelihood and its bias

Let's do some sanity check. Suppose that  $x_1, \dots, x_N$  are generated from a Gaussian distribution whose true parameters are  $\mu$  and  $\sigma^2$ . We expect the calculated parameters  $\mu_{ML}$  and  $\sigma_{ML}^2$  to be equal to  $\mu$  and  $\sigma^2$  respectively. Or put another way, we expect:

$$E(\mu_{ML}) = \mu$$

$$E(\sigma_{ML}^2) = \sigma^2$$

Is that true?

# Maximum likelihood and its bias

$$E(\mu_{ML}) = E\left(\frac{1}{N} \sum_{n=1}^N x_n\right) = \frac{1}{N} \sum_{n=1}^N E(x_n) = \mu$$

$$\begin{aligned} E(\mu_{ML}^2) &= E\left(\left(\frac{1}{N} \sum_{n=1}^N x_n\right)^2\right) \\ &= \frac{1}{N^2} \left( \sum_{1 \leq m \neq n \leq N} E(x_m x_n) + \sum_{n=1}^N E(x_n^2) \right) = \mu^2 + \frac{1}{N} \sigma^2 \end{aligned}$$

$$\begin{aligned} E(\sigma_{ML}^2) &= E\left(\frac{1}{N} \sum_{n=1}^N (x_n - \mu_{ML})^2\right) \\ &= \frac{1}{N} \sum_{n=1}^N E(x_n^2) - E(\mu_{ML}^2) = \frac{N-1}{N} \sigma^2 \end{aligned}$$

# Maximum likelihood and its bias

For a Gaussian distribution, the following estimate for the variance parameter is unbiased:

$$\tilde{\sigma}^2 = \frac{N}{N-1} \sigma_{ML}^2 = \frac{1}{N-1} \sum_{n=1}^N (x_n - \mu_{ML})^2$$



# Linear regression from a maximum likelihood perspective

## Problem

Assume that given the value of  $x_n$ , the corresponding value of  $t_n$  has a Gaussian distribution with a mean equal to the value  $y(x_n; w)$  and a variance  $\sigma^2$  (where the parameters  $w$  and  $\sigma^2$  are to be determined). Maximize the likelihood function:

$$p(t|x; w, \sigma^2) = \prod_{n=1}^N \mathcal{N}(t_n; y(x_n; w), \sigma^2)$$

# Linear regression from a maximum likelihood perspective

Again, we minimize the negative log function:

$$\begin{aligned} L &= -\log p(t|x; w, \sigma^2) \\ &= \frac{1}{2\sigma^2} \sum_{n=1}^N (y(x_n; w) - t_n)^2 + \frac{N}{2} \log \sigma^2 + \frac{N}{2} \log(2\pi) \end{aligned}$$

We see that maximizing the likelihood function for  $w$  is equivalent to minimizing the error function defined by:

$$E(w) = \frac{1}{2} \sum_{n=1}^N (y(x_n; w) - t_n)^2$$

# Linear regression from a maximum likelihood perspective

What has us gained from looking at the linear regression problem from a maximum likelihood perspective? Instead of a point estimate, we now have a predictive distribution:

$$p(\hat{t}|\hat{x}; w_{ML}, \sigma_{ML}^2) = \mathcal{N}(\hat{t}; y(\hat{x}; w_{ML}), \sigma_{ML}^2)$$

where

$$w_{ML} = (X^T X)^{-1} X^T t$$

$$\sigma_{ML}^2 = \frac{1}{N} \sum_{n=1}^N (y(x_n; w_{ML}) - t_n)^2$$

# Probability densities are integrand

When changing variable, we need to be aware that probability densities are integrand:

$$p(x)dx = p(g(y))dg(y) = p(g(y))g'(y)dy$$

For multivariate case:

$$p(x)dx = p(g(y)) \det \frac{\partial(x_1, \dots, x_N)}{\partial(y_1, \dots, y_N)} dy$$

# Transformation of densities

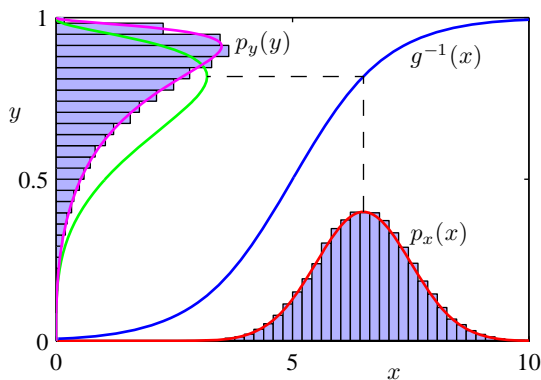
Consider the problem of finding the maximum for a probability density  $p(x)$ . Say the maximum happens when  $x = \hat{x}$ . Now we do a change of variable  $x = g(y)$ , does the maximum for the new probability density happens at  $\hat{y}$  where  $\hat{x} = g(\hat{y})$ ?

$$q(y) = p(g(y))g'(y)$$
$$q'(y) = p'(g(y))(g'(y))^2 + p(g(y))g''(y)$$

We see that this is usually not the case, unless  $g$  is a linear transformation.

# Transformation of densities

Figure: Transformation of the mode of a density



# Information

Intuitively, if we have two events  $x$  and  $y$  that are unrelated, the information gained from observing both of them should be the sum of the information gained from each of them separately:

$$h(x, y) = h(x) + h(y)$$

$$p(x, y) = p(x)p(y)$$

From this it's plausible to define  $h(x) = -\log_2 p(x)$ .

# Entropy

The entropy of a random variable  $x$  is defined as the expectation of the information  $h(x)$  with respect to the distribution  $p(x)$ :

$$H[x] = E(h) = \sum_x p(x)h(x) = - \sum_x p(x) \log_2 p(x)$$

When using logarithms to the base of 2, the units of  $H[x]$  are bits. From now on, we will switch to the use of natural logarithms in defining entropy, which is measured in units of *nats*.



# Maximum entropy for the discrete case

Let  $H(p) = -\sum_{n=1}^N p_i \log p_i$ , where  $0 \leq p_i \leq 1$ , it's easy to see that  $H(p)$  achieves its minimum 0 for unit vectors. When does  $H(p)$  achieves its maximum?

# Maximum entropy for the discrete case

Finding the maximum of  $H(p)$  under the constraint  $g(p) = \sum_{n=1}^N p_n - 1 = 0$  using Lagrange multiplier:

$$\begin{aligned}\nabla H(p) &= \lambda \nabla g(p) \\ -(\log p_n + 1) &= \lambda \\ p_n &= \frac{1}{N} \\ \max H(p) &= \log N\end{aligned}$$

# Differential entropy and its maximum

For the continuous case, we define the differential entropy to be:

$$H[x] = - \int p(x) \log p(x) dx$$

# Differential entropy and its maximum

Finding the maximum of  $H(p)$  under the following constraints:

$$\begin{aligned}\int_{-\infty}^{+\infty} p(x) dx &= 1 \\ \int_{-\infty}^{+\infty} xp(x) dx &= \mu \\ \int_{-\infty}^{+\infty} (x - \mu)^2 p(x) dx &= \sigma^2\end{aligned}$$

The maximum happens when  $p(x)$  is the Gaussian distribution:

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

and

$$\max H(p) = \frac{1}{2}(1 + \log(2\pi\sigma^2))$$

# Kullback-Leibler divergence

## Problem

Consider some unknown distribution  $p(x)$ . Suppose we have modelled  $p(x)$  using an approximating distribution  $q(x)$ . If we use  $q(x)$  to construct a coding scheme, what is the average additional amount of information required?

# Kullback-Leibler divergence

$$\begin{aligned} KL(p||q) &= - \int p(x) \log q(x) dx - (- \int p(x) \log p(x) dx) \\ &= - \int p(x) \log \frac{q(x)}{p(x)} dx \end{aligned}$$

This is also known as the relative entropy or Kullback-Leibler divergence, or KL divergence, between the distributions  $p(x)$  and  $q(x)$ .

# Kullback-Leibler divergence

If  $f$  is a convex function, then Jensen's inequality holds:

$$\begin{aligned} f(E(x)) &\leq E(f) \\ f\left(\sum_{n=1}^N p_n x_n\right) &\leq \sum_{n=1}^N p_n f(x_n) \\ f\left(\int x p(x) dx\right) &\leq \int p(x) f(x) dx \end{aligned}$$

Notice that  $-\log x$  is a convex function, we have:

$$KL(p||q) = \int p(x) \left(-\log \frac{q(x)}{p(x)}\right) dx \geq -\log \int p(x) \frac{q(x)}{p(x)} dx = 0$$

The equality will hold iff.  $q = p$ .

# Kullback-Leibler divergence

Minimizing the Kullback-Leibler divergence is equivalent to maximizing the likelihood function:

$$KL(p||q) \approx \frac{1}{N} \sum_{n=1}^N (-\log q(x_n; \theta) + \log p(x_n))$$

The first term is the negative log likelihood function for  $\theta$  under the distribution  $q(x; \theta)$  evaluated using the training set.



# Conditional entropy

On average, if value for one random variable is already known, what is the additional information needed to specify value for another random variable?

$$H[y|x] = - \iint p(x, y) \log p(y|x) dx dy$$
$$H[x, y] = H[y|x] + H[x]$$

# Mutual information

For two random variables, are they "close" to being independent?

$$\begin{aligned} I[x, y] &= KL(p(x, y) || p(x)p(y)) \\ &= - \iint p(x, y) \log \frac{p(x)p(y)}{p(x, y)} dx dy \end{aligned}$$

It's easy to see that:

$$I[x, y] = H[x] - H[x|y] = H[y] - H[y|x]$$

# Model parameters

Denote the training data set by  $\mathcal{D}$ , and the parameters in the model by  $w$ .

- $p(w)$  is our assumptions about  $w$  before observing  $\mathcal{D}$ .
- $p(\mathcal{D}|w)$  is the likelihood function.
- $p(w|\mathcal{D})$  is the uncertainty in  $w$  after we have observed  $\mathcal{D}$ .

We have:

$$\begin{aligned} p(w|\mathcal{D}) &= \frac{p(\mathcal{D}|w)p(w)}{p(\mathcal{D})} \\ &= \frac{p(\mathcal{D}|w)p(w)}{\int p(\mathcal{D}|w)p(w)dw} \end{aligned}$$

# Regularization

When choosing the model parameters  $w$ , instead of maximizing the likelihood function  $p(\mathcal{D}|w)$ , we maximize the posterior probability  $p(w|\mathcal{D})$ :

$$-\log p(w|\mathcal{D}) = -\log p(\mathcal{D}|w) - \log p(w) + \log p(\mathcal{D})$$

Say each  $w_m$  conforms to a Gaussian distribution:

$$p(w) = p(w; \sigma^2) = \prod_{m=0}^M \mathcal{N}(w_m; 0, \sigma^2)$$

Then we have:

$$-\log p(w|\mathcal{D}) = -\log p(\mathcal{D}|w) + \frac{1}{2\sigma^2} \sum_{m=0}^M w_m^2 + \text{const}$$

The second term on the right hand side is indeed the penalty term.

# Bayesian machine learning

If we are interested in the distribution of  $t$  given both  $x$  and  $\mathcal{D}$ , taking into consideration the uncertainty in the value of  $w$ , we have the fully Bayesian treatment:

$$p(t|x, \mathcal{D}) = \int p(t|x, w)p(w|\mathcal{D})dw$$

- The fully Bayesian treatment averages over all possible models:
  - Less likely to lead to over-fitting.
  - Prefer models of intermediate complexity.
- Integrating over the space of parameters is typically infeasible.