Machine Learning CMPT 726

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2021-10-25

Probabilistic Interpretation of Linear Regression

Probabilistic Model

Suppose we know how the data was generated (known as the "data generating process"):

$$y = \overrightarrow{w}^{\mathsf{T}} \overrightarrow{x} + \sigma \epsilon$$
 ,where $\epsilon \sim \mathcal{N}(0,1)$

 \overrightarrow{w} and σ are unknown parameters

We observe many tuples (\vec{x}, y) , denoted as $\{(\vec{x}_i, y_i)\}_{i=1}^N =: \mathcal{D}$, which we assume are generated i.i.d. from the process above.

Terminology: i.i.d. stands for "independent and identically distributed", which means that each tuple (\vec{x}_i, y_i) is independent of other tuples and has the same joint distribution as any other tuple.

Goal: Estimate the values of unknown parameters from observations (known as "parameter estimation").

Probabilistic Model

A probabilistic model assigns a probability to every possible observation.

We choose a probabilistic model based on the data generating process:

Recall: data generating process: $y=\vec{w}^{\mathsf{T}}\vec{x}+\sigma\epsilon$,where $\epsilon\sim\mathcal{N}(0,1)$

Recall: if
$$Z \sim \mathcal{N}(0,1)$$
, $\mu + \sigma Z \sim \mathcal{N}(\mu, \sigma^2)$

So,
$$y | \vec{x}, \vec{w}, \sigma \sim \mathcal{N}(\vec{w}^{\mathsf{T}} \vec{x}, \sigma^2)$$

We can write down the expression for $p(y|\vec{x}, \vec{w}, \sigma)$:

Recall: If
$$X \sim \mathcal{N}(\mu, \sigma^2)$$
, $p(x) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right)$
$$p(y|\vec{x}, \vec{w}, \sigma) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(y-\vec{w}^\top \vec{x})^2}{2\sigma^2}\right)$$

Parameter Estimation

We have two unknown parameters: \overrightarrow{w} and σ .

For the purposes of making predictions, we care mostly about \overrightarrow{w} and so will focus on estimating \overrightarrow{w} .

In general:

The parameters we care about are known as parameters of interest.

The parameters we don't care about are known as nuisance parameters.

In this case, \vec{w} is the parameter of interest, and σ is the nuisance parameter.

Idea: Find the parameter value at which the probability of observing the data is maximized.

Step 1: Derive the joint probability density of the observations $p(\mathcal{D}|\vec{w}, \sigma)$.

$$p(\mathcal{D}|\vec{w},\sigma):=p(y_1,\ldots,y_N|\vec{x}_1,\ldots,\vec{x}_N,\vec{w},\sigma)$$

$$=\prod_{i=1}^N p(y_i|\vec{x}_1,...,\vec{x}_N,\vec{w},\sigma)$$
 (Conditioned on $\vec{x}_1,...,\vec{x}_N$, the y_i 's are independent)

$$=\prod_{i=1}^{N}p(y_i|\vec{x}_i,\vec{w},\sigma)$$
 (Conditioned on each \vec{x}_i , y_i and \vec{x}_j 's for $j \neq i$ are independent)

$$= \prod_{i=1}^{N} \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(y_i - \vec{w}^{\mathsf{T}} \vec{x}_i)^2}{2\sigma^2}\right)$$

$$:= \mathcal{L}(\overrightarrow{w}, \sigma; \{(\overrightarrow{x}_i, y_i)\}_{i=1}^N) = \mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D})$$

"Likelihood function"

Step 2: Find the value of the parameter of interest that maximizes the likelihood function.

$$\widehat{\overrightarrow{w}}_{\text{MLE}} = \arg \max_{\overrightarrow{w}} \mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D})$$

To find it, we can try computing the gradient and set it to zero.

$$\nabla_{\vec{w}} \mathcal{L}(\vec{w}, \sigma; \mathcal{D}) = \nabla_{\vec{w}} \left(\prod_{i=1}^{N} p(y_i | \vec{x}_i, \vec{w}, \sigma) \right)$$

$$= \sum_{i=1}^{N} \left(\nabla_{\vec{w}} p(y_i | \vec{x}_i, \vec{w}, \sigma) \prod_{j \neq i} p(y_j | \vec{x}_j, \vec{w}, \sigma) \right)$$
Unwie

Step 2: Find the value of the parameter of interest that maximizes the likelihood function.

$$\widehat{\overrightarrow{w}}_{\text{MLE}} = \arg \max_{\overrightarrow{w}} \mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D})$$

When the base of the log is not shown, it defaults to e (natural number)

Instead use the fact that $\widehat{\overrightarrow{w}}_{\text{MLE}} = \arg\max_{\overrightarrow{w}} \mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D}) = \arg\max_{\overrightarrow{w}} \log\mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D}).$

)This is true because $x \mapsto \log(x)$ is strictly increasing, i.e.: as x increases, $\log(x)$ increases)

$$\log \mathcal{L}(\vec{w}, \sigma; \mathcal{D}) = \log \prod_{i=1}^{N} p(y_i | \vec{x}_i, \vec{w}, \sigma) = \sum_{i=1}^{N} \log p(y_i | \vec{x}_i, \vec{w}, \sigma)$$
 "Log-likelihood function"

-1 0 1 2 3 4

This turns the expression into a sum, which is easy to differentiate.

$$\nabla_{\overrightarrow{w}} \log \mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D}) = \nabla_{\overrightarrow{w}} \left(\sum_{i=1}^{N} \log p(y_i | \overrightarrow{x}_i, \overrightarrow{w}, \sigma) \right) = \sum_{i=1}^{N} \nabla_{\overrightarrow{w}} \log p(y_i | \overrightarrow{x}_i, \overrightarrow{w}, \sigma)$$

Step 2: Find the value of the parameter of interest that maximizes the likelihood function.

$$\widehat{\vec{w}}_{\text{MLE}} = \arg \max_{\vec{w}} \mathcal{L}(\vec{w}, \sigma; \mathcal{D}) = \arg \max_{\vec{w}} \log \mathcal{L}(\vec{w}, \sigma; \mathcal{D})$$

$$\log \mathcal{L}(\vec{w}, \sigma; \mathcal{D}) = \sum_{i=1}^{N} \log p(y_i | \vec{x}_i, \vec{w}, \sigma)$$

$$= \sum_{i=1}^{N} \log \left(\frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(y_i - \vec{w}^\top \vec{x}_i)^2}{2\sigma^2} \right) \right)$$

$$= \sum_{i=1}^{N} \left(-\frac{(y_i - \vec{w}^\top \vec{x}_i)^2}{2\sigma^2} - \log\sqrt{2\pi\sigma^2} \right)$$

$$= -\sum_{i=1}^{N} \frac{(y_i - \vec{w}^\top \vec{x}_i)^2}{2\sigma^2} - N\log\sqrt{2\pi\sigma^2}$$

$$= -\frac{1}{2\sigma^2} \sum_{i=1}^{N} (y_i - \vec{w}^\top \vec{x}_i)^2 - N\log\sqrt{2\pi\sigma^2}$$

$$\widehat{\overrightarrow{w}}_{\text{MLE}} = \operatorname{argmax} \log \mathcal{L}(\overrightarrow{w}, \sigma; \mathcal{D})$$

$$= \operatorname{arg} \max_{\overrightarrow{w}} \left(-\frac{1}{2\sigma^2} \sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\mathsf{T}} \overrightarrow{x}_i)^2 - N \log \sqrt{2\pi\sigma^2} \right)$$

$$= \operatorname{arg} \min_{\overrightarrow{w}} \left(\frac{1}{2\sigma^2} \sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\mathsf{T}} \overrightarrow{x}_i)^2 + N \log \sqrt{2\pi\sigma^2} \right)$$

$$= \operatorname{arg} \min_{\overrightarrow{w}} \frac{1}{2\sigma^2} \sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\mathsf{T}} \overrightarrow{x}_i)^2$$

$$= \operatorname{arg} \min_{\overrightarrow{w}} \sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\mathsf{T}} \overrightarrow{x}_i)^2$$
Compare to

Compare to the loss function of OLS:

$$L(\vec{w}) = \sum_{i=1}^{N} (y_i - \vec{w}^{\mathsf{T}} \vec{x}_i)^2$$

Frequentist vs. Bayesian Statistics

Previously we treated parameters as fixed (albeit unknown) quantities.

Suppose we know what values of the parameters are more plausible compared to others. This is known as **prior belief** or **prior knowledge**.

We can represent this prior belief as a probability distribution over parameters, which is known as a **prior probability distribution**.

Instead of treating parameters as fixed quantities, we treat the parameters as random variables that follow the prior probability distribution.

This latter approach is known as **Bayesian statistics**, whereas the former approach is known as **frequentist statistics**.

In addition to the probabilistic model, we have a prior distribution $p(\vec{\theta})$ over the parameters of interest $\vec{\theta}$.

Idea: Find the parameter value that maximizes the conditional probability over parameters given the observations $p\left(\vec{\theta} \middle| \mathcal{D}\right)$.

We can use Bayes' rule to find $p\left(\overrightarrow{\theta}\middle|\mathcal{D}\right)$:

"Prior" "Likelihood"

"Posterior"

$$p\left(\vec{\theta}\middle|\mathcal{D}\right) = \frac{p\left(\vec{\theta}\right)p(\mathcal{D}\middle|\vec{\theta}\right)}{p(\mathcal{D})}$$

$$\hat{\vec{\theta}}_{MAP} = \arg\max_{\vec{\theta}} p\left(\vec{\theta}\middle|\mathcal{D}\right)$$

The term "likelihood" is used in broader contexts than "likelihood function". While its meaning coincides with "likelihood function" in this context, it could mean the probability of a single observation $p(y|\vec{x},\vec{\theta})$ or any $p(\vec{x}|\vec{y})$ where we observe \vec{x} and would like to infer an unobserved \vec{y} based on \vec{x} .

Consider the same probabilistic model as before:

$$y|\vec{x}, \vec{w}, \sigma \sim \mathcal{N}(\vec{w}^{\mathsf{T}} \vec{x}, \sigma^{2})$$

$$p(y|\vec{x}, \vec{w}, \sigma) = \frac{1}{\sqrt{2\pi\sigma^{2}}} \exp\left(-\frac{(y - \vec{w}^{\mathsf{T}} \vec{x})^{2}}{2\sigma^{2}}\right)$$

We choose the following prior distribution over the parameter of interest \vec{w} :

$$\overrightarrow{w}|\sigma \sim \mathcal{N}\left(\overrightarrow{0}, \frac{\sigma^2}{\lambda}I\right)$$

$$p(\overrightarrow{w}|\sigma) = \frac{1}{\sqrt{(2\pi)^n \det\left(\frac{\sigma^2}{\lambda}I\right)}} \exp\left(-\frac{1}{2}\overrightarrow{w}^{\mathsf{T}}\left(\frac{\sigma^2}{\lambda}I\right)^{-1}\overrightarrow{w}\right) = \frac{1}{\sqrt{\left(\frac{2\pi\sigma^2}{\lambda}\right)^n}} \exp\left(-\frac{\lambda}{2\sigma^2}||\overrightarrow{w}||_2^2\right)$$

$$\begin{split} &\widehat{\overrightarrow{w}}_{\text{MAP}} = \arg\max_{\overrightarrow{w}} p(\overrightarrow{w}|\mathcal{D}, \sigma) \\ &= \arg\max_{\overrightarrow{w}} \log p(\overrightarrow{w}|\mathcal{D}, \sigma) \\ &= \arg\max_{\overrightarrow{w}} \log \left(\frac{p(\overrightarrow{w}|\sigma)p(\mathcal{D}|\overrightarrow{w}, \sigma)}{p(\mathcal{D}|\sigma)} \right) \\ &= \arg\max_{\overrightarrow{w}} [\log p(\overrightarrow{w}|\sigma) + \log p(\mathcal{D}|\overrightarrow{w}, \sigma) - \log p(\mathcal{D}|\sigma)] \\ &= \arg\max_{\overrightarrow{w}} [\log p(\overrightarrow{w}|\sigma) + \log p(\mathcal{D}|\overrightarrow{w}, \sigma)] \\ &= \arg\max_{\overrightarrow{w}} [\log p(\overrightarrow{w}|\sigma) + \log p(y_1, \dots, y_N|\overrightarrow{x}_1, \dots, \overrightarrow{x}_N, \overrightarrow{w}, \sigma)] \\ &= \arg\max_{\overrightarrow{w}} [\log p(\overrightarrow{w}|\sigma) + \log \prod_{i=1}^{N} p(y_i|\overrightarrow{x}_i, \overrightarrow{w}, \sigma)] \\ &= \arg\max_{\overrightarrow{w}} [\log p(\overrightarrow{w}|\sigma) + \sum_{i=1}^{N} \log p(y_i|\overrightarrow{x}_i, \overrightarrow{w}, \sigma)] \end{split}$$

$$\begin{split} \widehat{\vec{w}}_{\text{MAP}} &= \arg\max_{\vec{w}} [\log p(\vec{w}|\sigma) + \sum_{i=1}^{N} \log p(y_{i}|\vec{x}_{i}, \vec{w}, \sigma)] \\ &= \arg\max_{\vec{w}} \left[\log \left(\frac{1}{\sqrt{\left(\frac{2\pi\sigma^{2}}{\lambda} \right)^{n}}} \exp\left(-\frac{\lambda}{2\sigma^{2}} ||\vec{w}||_{2}^{2} \right) \right) + \sum_{i=1}^{N} \log \left(\frac{1}{\sqrt{2\pi\sigma^{2}}} \exp\left(-\frac{(y_{i} - \vec{w}^{\top} \vec{x}_{i})^{2}}{2\sigma^{2}} \right) \right) \right] \\ &= \arg\max_{\vec{w}} \left[-\frac{\lambda}{2\sigma^{2}} ||\vec{w}||_{2}^{2} - \log \left(\sqrt{\left(\frac{2\pi\sigma^{2}}{\lambda} \right)^{n}} \right) + \sum_{i=1}^{N} \left(-\frac{(y_{i} - \vec{w}^{\top} \vec{x}_{i})^{2}}{2\sigma^{2}} - \log\sqrt{2\pi\sigma^{2}} \right) \right] \\ &= \arg\max_{\vec{w}} \left[-\frac{\lambda}{2\sigma^{2}} ||\vec{w}||_{2}^{2} + \sum_{i=1}^{N} \left(-\frac{(y_{i} - \vec{w}^{\top} \vec{x}_{i})^{2}}{2\sigma^{2}} \right) \right] \end{split}$$

$$\widehat{\vec{w}}_{MAP} = \arg \max_{\vec{w}} \left[-\frac{\lambda}{2\sigma^2} \|\vec{w}\|_2^2 + \sum_{i=1}^N \left(-\frac{(y_i - \vec{w}^\top \vec{x}_i)^2}{2\sigma^2} \right) \right]$$

$$= \arg \max_{\vec{w}} -\frac{1}{2\sigma^2} \left(\lambda \|\vec{w}\|_2^2 + \sum_{i=1}^N (y_i - \vec{w}^\top \vec{x}_i)^2 \right)$$

$$= \arg \min_{\vec{w}} \frac{1}{2\sigma^2} \left(\lambda \|\vec{w}\|_2^2 + \sum_{i=1}^N (y_i - \vec{w}^\top \vec{x}_i)^2 \right)$$

$$= \arg \min_{\vec{w}} \left(\lambda \|\vec{w}\|_2^2 + \sum_{i=1}^N (y_i - \vec{w}^\top \vec{x}_i)^2 \right)$$

Compare to the loss function of ridge regression: $L(\vec{w}) = \sum_{i=1}^{r} (y_i - \vec{w}^{\mathsf{T}} \vec{x}_i)^2 + \lambda ||\vec{w}||_2^2$

OLS vs. Ridge Regression (Deterministic Interpretation)

Model: $\vec{y} = \vec{w}^{\mathsf{T}} \vec{x}$

Parameters: \overrightarrow{w}

OLS:

Loss function:

$$L(\overrightarrow{w}) = \sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\mathsf{T}} \overrightarrow{x}_i)^2$$
$$= \|\overrightarrow{y} - X\overrightarrow{w}\|_2^2$$

Optimal Parameters:

$$\overrightarrow{w}^* := \arg\min_{\overrightarrow{w}} L(\overrightarrow{w}) = (X^T X)^{-1} X^T \overrightarrow{y}$$

Ridge Regression:

Loss function:

$$L(\vec{w}) = \sum_{i=1}^{N} (y_i - \vec{w}^{\mathsf{T}} \vec{x}_i)^2 + \lambda ||\vec{w}||_2^2$$
$$= ||\vec{y} - X\vec{w}||_2^2 + \lambda ||\vec{w}||_2^2$$

Optimal Parameters:

$$\vec{w}^* := \arg\min_{\vec{w}} L(\vec{w}) = (X^T X + \lambda I)^{-1} X^T \vec{y}$$

OLS vs. Ridge Regression (Probabilistic Interpretation)

Probabilistic Model: $y|\vec{x}, \vec{w}, \sigma \sim \mathcal{N}(\vec{w}^{\top}\vec{x}, \sigma^2)$; Prior: $\vec{w}|\sigma \sim \mathcal{N}\left(\vec{0}, \frac{\sigma^2}{\lambda}I\right)$

Parameters: \overrightarrow{w}

Nuisance Parameter: σ

OLS:

MLE estimate:

$$\widehat{\overrightarrow{w}}_{\text{MLE}} = \arg\min_{\overrightarrow{w}} \sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\top} \overrightarrow{x}_i)^2$$

$$= \arg\min_{\overrightarrow{w}} ||\overrightarrow{y} - X\overrightarrow{w}||_2^2$$

$$= (X^{\top} X)^{-1} X^{\top} \overrightarrow{y}$$

Ridge Regression:

MAP estimate:

$$\widehat{\overrightarrow{w}}_{\text{MAP}} = \arg\min_{\overrightarrow{w}} \left[\sum_{i=1}^{N} (y_i - \overrightarrow{w}^{\top} \overrightarrow{x}_i)^2 + \lambda ||\overrightarrow{w}||_2^2 \right]$$

$$= \arg\min_{\overrightarrow{w}} [||\overrightarrow{y} - X\overrightarrow{w}||_2^2 + \lambda ||\overrightarrow{w}||_2^2]$$

$$= (X^{\top} X + \lambda I)^{-1} X^{\top} \overrightarrow{y}$$

Takeaways

There are deep connections between deterministic and probabilistic formulations of machine learning methods.

Many methods have both deterministic and probabilistic interpretations.

Primary loss function is related to the likelihood.

Regularizer is related to the prior.

Square losses are related to (univariate) Gaussian observation noise.

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

Squared L2 losses are related to (multivariate) isotropic Gaussian observation noise.

$$p(\vec{x}) = \frac{1}{\sqrt{(2\pi\sigma^2)^n}} \exp\left(-\frac{1}{2\sigma^2} ||\vec{x} - \vec{\mu}||_2^2\right)$$

(Full) Bayesian Inference

Bayesian inference refers to computing the full posterior:

$$p(\vec{\theta}|\mathcal{D}) = \frac{p(\vec{\theta})p(\mathcal{D}|\vec{\theta})}{p(\mathcal{D})}$$

Compare to MAP:

$$\hat{\vec{\theta}}_{MAP} = \arg \max_{\vec{\theta}} p(\vec{\theta}|\mathcal{D})$$

Unlike MAP estimation, it does not just produce a single value for the parameter estimate (known as a **point estimate**). Instead, it produces a full distribution over possible parameter values.

Can be extremely computationally challenging - computing the posterior exactly is intractable for all but some special cases. Must typically rely on sampling or approximations.

Note on the Term "Inference"

Parameter estimation is also known as **inference** (especially in the context of Bayesian inference), since the purpose is to infer the value of unknown parameters from observed data.

Note: In machine learning, the term "inference" is overloaded and can sometimes mean testing/making predictions (as opposed to training).

This can get confusing, because parameter estimation in a probabilistic model corresponds to training the model, which is also known as inference. On the other hand, testing the model is also sometimes known as inference. So, inference can mean either training or testing!

Meaning depends on context:

- "inference time", "training and inference": means testing
- "Bayesian inference", "probabilistic inference", "parameter inference": means training
- "inference procedure": ambiguous

In this course, we will avoid using the term as much as possible to minimize confusion.

Quiz Practice

Q1: Which of the following statements about the difference between MLE and MAP parameter estimates is NOT true?

- (A) MLE does not depend on the prior, whereas MAP does
- (B) MLE depends on the likelihood function, whereas MAP does not
- (C) The derivation of MLE does not use Bayes' rule, whereas the derivation of MAP does
- (D) Ridge regression can be interpreted as MAP, but not as MLE
- (E) MAP corresponds to optimizing a loss function with a regularizer on the parameters, whereas MLE corresponds to optimizing a loss function without a regularizer
- (F) All are true