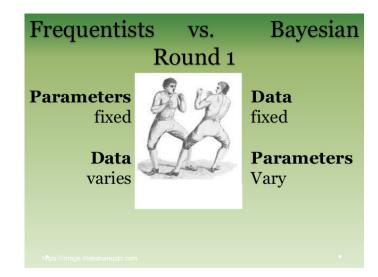


Overview

- Model fitting
- Model selection

Frequentist vs Bayesian

- Probabilistic Model p(y, x, w)
 - Frequentists: w is a parameter that should be estimated by model fitting
 - **Bayesians**: w is a random variable that has a prior distribution p(w)
 - How to set p(w)??



Example: Linear regression, what are parameters here?

$$y \sim w_0 + \boldsymbol{w}\boldsymbol{x} + e, e \sim N(0, \sigma^2)$$

 $y \sim N(w_0 + \boldsymbol{w}\boldsymbol{x}, \sigma^2)$

An estimator

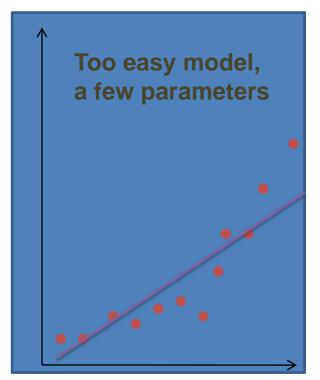
- $\hat{w} = \delta(D)$ (some function of your data) an estimator
- Optimal parameter values? → there can be many ways to compute them (MLE, shrinkage...)
 - Compare Bayesian: given estimators w^1 and w^2 , we can compare them! $p(w^1|D) > p(w^2|D)$
 - There is no easy way to compare estimators in frequentist tradition

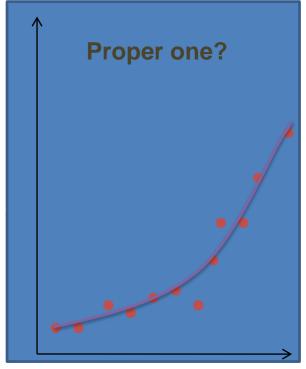
Example: Linear regression

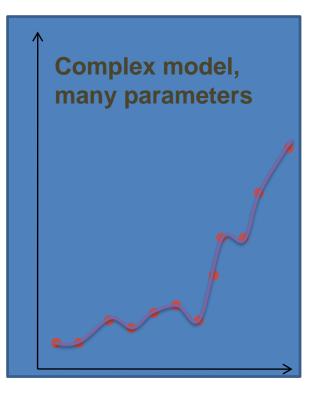
- Estimator 1: $\mathbf{w} = (X^T X)^{-1} X^T Y$ (maximum likelihood)
- Estimator 2: w = (0, ..., 0, 1)
- Which one is better?
 - A comparison strategy is needed!

Overfitting

Complex model can overfit your data







Overfitting: solutions

Observed: Maximum likelihood can lead to overfitting.

Solutions

- Selecting proper parameter values
 - Regularized risk minimization
- Selecting proper model type, for ex. number of parameters
 - Houldout method
 - Cross-validation

Model selection

- Given a model, choose the optimal parameter values
 - Decision theory
- Define loss $L(Y, \hat{y})$
 - How much we loose in guessing true Y incorrectly
- If we know the true distribution p(y, x|w) then we choose \hat{y}

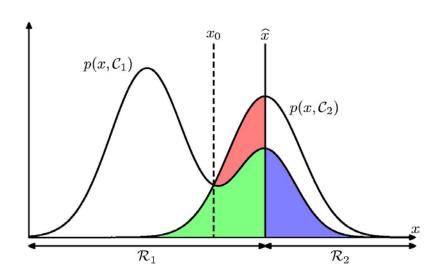
$$\min_{\hat{f}} EL(y, \hat{y}) = \min_{\hat{y}} \int L(y, \hat{y}) p(y, x|w) dx dy$$

Model selection

Example: Spam classification

Loss for incorrect classifying mails and spams

$$-L_{12}=1, L_{21}=100$$



Loss functions

- How to define loss function?
 - No unique choice, often defined by application
 - Normal practice: Choose the loss related to minus loglikelihood

Example: Predicting the amount of the product at the storage:

$$L(Y, \hat{y}) = \begin{cases} 10, +\hat{y}/Y \ \hat{y} \ge Y \\ 1000, \hat{y} < Y \end{cases}$$

Example: Compute loss function related to

Normal distribution

Guess why such loss function was chosen

Loss functions

Classification problems

- Common loss function
$$L(Y, \hat{y}) = \begin{cases} 1, Y = \hat{y} \\ 0, Y \neq \hat{y} \end{cases}$$

When minimizing the loss, equivalent to misclassification rate

Model selection

- Problem: true model and true w are unknown → can not compute expected loss!
- How to find an optimal model?
 - Consider what expected loss (**risk**) depends on $R(Y, \hat{y}) = E[L(Y, \hat{y}(X, D))]$
- Random factors:
 - − D − training set
 - Y, X data to be predicted (validation set)

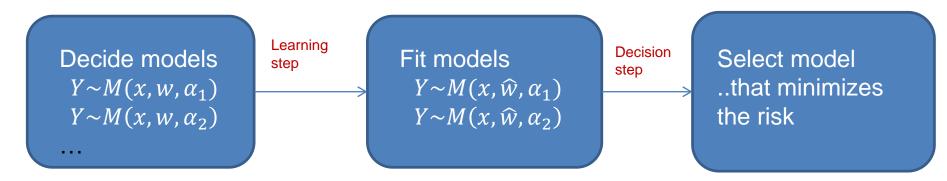
- Simplify the risk estimation:
 - Fix D as a particular training set T
 - Fix Y,X as a particular validation set V
- Risk becomes (empirical risk)

$$\widehat{R}(y,\widehat{y}) = \frac{1}{|V|} \sum_{(X,Y) \in V} L(Y,\widehat{y}(X,T))$$

- Estimator is fit by Maximum Likelihood using training set
- Risk estimated by using validation set
- Model with minimum empirical risk is selected

General model selection strategy

• Given data $D = \{X_i, Y_i, i = 1 ... n\}$



- When fitting data, Maximum Likelihood is usually used
- α_i can be different things:
 - Type of distribution
 - Number of variables in the model
 - Regulatization parameter value
 - ...

Divide into training, validation and test sets

Training Validation Test

Choose proportions in some way

 Given: training, validation, test sets and models to select between

M1(?,?)

M2(?,?,?)

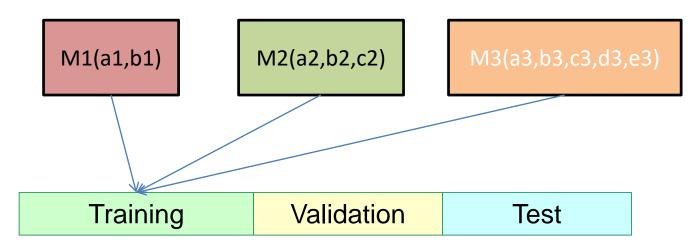
M3(?,?,?,?,?)

Training

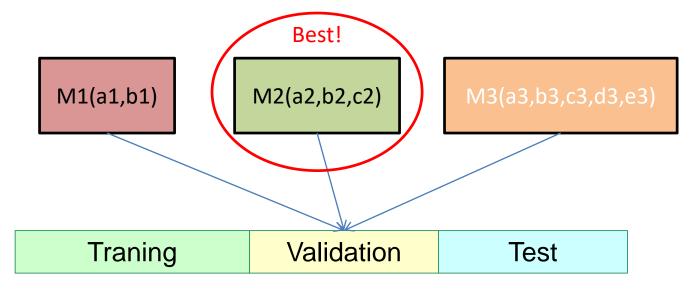
Validation

Test

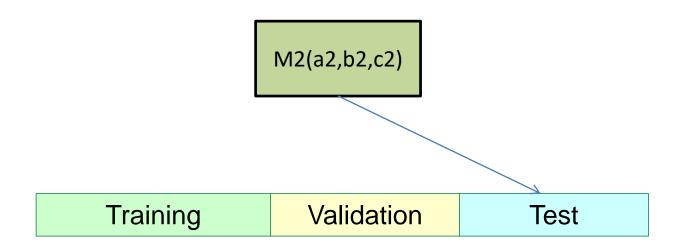
 Training set is to used for fitting models to the dataset by using maximum likelihood

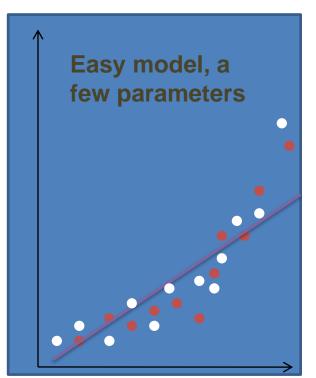


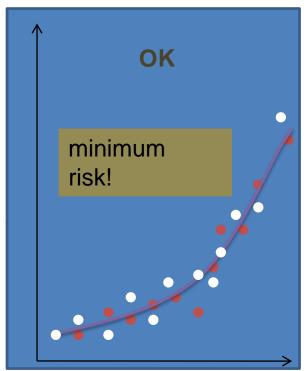
 Validation set is used to choose the best model (lowest risk)

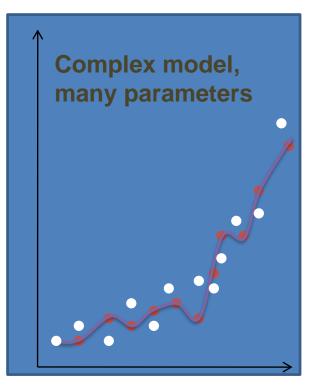


 Test set is used to test a performance on a new data









Holdout in R

- How to partition into train/test?
 - Use set.seed(12345) in the labs to get identical results

```
n=dim(data)[1]
set.seed(12345)
id=sample(1:n, floor(n*0.7))
train=data[id,]
test=data[-id,]
```

How to partition into train/valid/test?

```
n=dim(data)[1]
set.seed(12345)
id=sample(1:n, floor(n*0.4))
train=data[id,]

id1=setdiff(1:n, id)
set.seed(12345)
id2=sample(id1, floor(n*0.3))
valid=data[id2,]

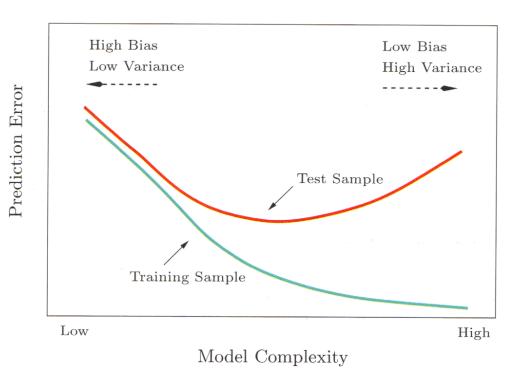
id3=setdiff(id1,id2)
test=data[id3,]
```

Bias-variance tradeoff

- Bias of an estimator $Bias(\hat{y}(x_0)) = E[\hat{y}(x_0) f(x_0)], f(x_0)$ is expected response
 - If $Bias(\hat{y}(x_0)) = 0$, the estimator is **unbiased**
 - ML estimators are asymptotically unbiased if the model is enough complex
 - However, unbiasedness does not mean a good choice!

Bias-variance tradeoff

• Assume loss is $L(Y, \hat{y}) = (Y - \hat{y})^2$ $R(Y(x_0), \hat{y}(x_0)) = \sigma^2 + Bias^2(\hat{y}(x_0)) + Var(\hat{y}(x_0))$



When loss is not quadratic, no such nice formula exist

Cross-validation

- Compared to holdout method:
 - Why do we use only some portion of data for training- can we use more (increase accuracy)?

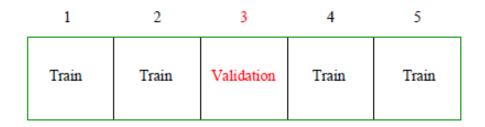
Cross-validation (Estimates Err)

K-fold cross-validation (rough scheme, show picture):

- 1. Permute the observations randomly
- 2. Divide data-set in K roughly equally-sized subsets
- 3. Remove subset #i and fit the model using remaining data.
- 4. Predict the function values for subset #i using the fitted model.
- 5. Repeat steps 3-4 for different i
- 6. CV= squared difference between observed values and predicted values (another function is possible)

Cross-validation

Cross-validation



Note: if K=N then method is *leave-one-out* crossvalidation.

$$\kappa: \{1, \ldots, N\} \mapsto \{1, \ldots, K\}$$

K-fold cross-validation:
$$CV =$$

$$\frac{1}{N}\sum_{i=1}^{N}L(Y_{i},\hat{y}^{-k(i)}(x_{i}))$$

What to do if N is not a multiple of K?

Cross-validation vs Holdout

Holdout is easy to do (a few model fits to each data)

 Cross validation is computationally demanding (many model fits)

- Holdout is applicable for large data
 - Otherwise, model selection performs poorly
- Cross validation is more suitable for smaller data

Analytical methods

- Analytical expressions to select models
 - AIC (Akaike's information criterion)

Idea: Instead of $R(Y, \hat{y}) = E[L(Y, \hat{y}(X, D))]$ consider **in-sample** risk (only Y in D is random):

$$R_{in}(Y, \hat{y}) = \frac{1}{N} \sum_{i=1}^{N} E_{Y_i} [L(Y_i, \hat{y}(X, D)) | D, X \in D]$$

Analytical methods

One can show that

$$R_{in}(Y,\hat{y}) \approx R_{train} + \frac{2}{N} \sum_{i} cov(\hat{y}_i, Y_i)$$
 where $R_{train} = \sum_{X_i, Y_i \in T} L(Y_i, \hat{y}_i)$

- Recall, degrees of freedom df(model) = $\frac{1}{\sigma^2} \sum_i cov(\hat{y}_i, Y_i)$
 - When model is linear, df is the number of parameters.
- If loss is defined by minus two loglikelihood, $AIC \equiv -2loglik(D) + 2df(model)$

Model selection

Example Computer Hardware Data Set: performance measured for various processors and also

- Cycle time
- Memory
- Channels
- ...

Build model predicting performance



Cross-validatation

Try models with different predictor sets

```
data=read.csv("machine.csv", header=F)
library(cvTools)
```

```
fit1=lm(V9~V3+V4+V5+V6+V7+V8, data=data)
fit2=lm(V9~V3+V4+V5+V6+V7, data=data)
fit3=lm(V9~V3+V4+V5+V6, data=data)
f1=cvFit(fit1, y=data$V9, data=data,K=10,
foldType="consecutive")
f2=cvFit(fit2, y=data$V9, data=data,K=10,
foldType="consecutive")
f3=cvFit(fit3, y=data$V9, data=data,K=10,
foldType="consecutive")
res=cvSelect(f1,f2,f3)
plot(res)
```

