Resampling in Statistics

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If you feel sleepy...



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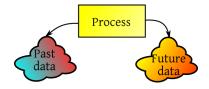
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Why do we hope to be better prepared for the future by analyzing past data?

Because we believe that we are going to face similar data again in the future.



An analogy

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If the syllabus changes, the interest in solving the past years' papers dwindles.

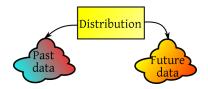
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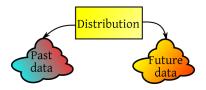
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There are other types of data generations processes (e.g., time series or spatial data). But for resampling we work almost exclusively with IID data.

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In step 0, we shall take just the opposite view point. We don't have any data, just the specification of a random experiment. A complete specification. No unknown parameters or anything. Can we generate the data then? The answer is yes, we can, by just running the random experiment.

Enter computers

Generating data from a specified random experiment is called simulation or Monte Carlo techniques.

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```
sample(6,100,replace=TRUE)
rnorm(100,mean=3,sd=0.5) rpois(100,lambda=3)
rbinom(100,size=10,prob=0.2)
rweibull(100,shape=3,scale=1)
```

Step 1: Quest for future data

We start with a data set, an IID one. So we know that there is some random experiment up there in the clouds.

We have done something (estimation, prediction etc) to prepare for the future.

We want to test our preparation.

Step 1: Quest for future data

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We may just wait until future data arrive, and then check. But it's like testing for a poison by actually swallowing it!

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We want some pretty good approximation of the future data. And that's exactly what resampling technique is all about:

We have past data, which are IID from some distribution with is not completely known. We want an approximation of future data from the same (unknown) distribution.

Three major approaches

- Cross-validation: the simplest of them all. Thanks to the popularity of deep learning, this old technique has now gained a lot of currency.
- Permutation test: A more sophisticated approach, but somewhat limited in scope.
- Bootstrap: The master technique that has a nice theory behind it, and is quite general.

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How can he use this?

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Reserve a subset for the exam, work out the rest.

Cross-validation

Split past data into two parts, a big part as the training data, and the rest as the test part. Fit all your models etc on the training part only, and then test out the performance on the test set!

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Implementation details may vary.

An example

There is a famous data set called the iris data about 150 iris flowers:

```
> head(iris)
Sepal.Length Sepal.Width Petal.Length Petal.Width Species
1 5.1 3.5 1.4 0.2 setosa
2 4.9 3.0 1.4 0.2 setosa
...
149 6.2 3.4 5.4 2.3 virginica
150 5.9 3.0 5.1 1.8 virginica
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Suppose that we want to come up with a rule by which you can identify the species based on these measurements.

Many reasonable techniques possible. Which is the best?

Different techniques

One possible method: Use only sepal lengths. Find the average sepal length for each species. When you are given a new flower, measure its sepal length, and put in the species that has the closest mean.

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```
useSL = function(pastData, newInput) {
ind.setosa = pastData[,5] == "setosa"
mean.setosa = mean(pastData[ind.setosa,1])
...
meanVector = c(mean.setosa,mean.virginica,mean.versicolor)
for(i in 1:nrow(newInput))
predictedClass[i] =
which.min(abs(newInput[i,1]-meanVector))

c('setosa','virginica','versicolor')[predictedClass]
}
```

But why not use...

- petal length?
- ► all the variables?
- some weighted average?
- median instead of mean?
- some statistical model?

We had 50 flowers of each species. Let's set aside 10 of each species, and try out your methods on the remaining 40*3 = 120 flowers.

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Now, we actually know the true species of these flowers. So we can readily find out the number of misclassifications that each method leads to.

The one leading to least number of misclassifications may be considered the best.

R code

```
mixup =
c(sample(50,50),50+sample(50,50),100+sample(50,50))
train = mixup[c(1:30,51:80,101:130)]

pred1=useSL(iris[train,],iris[-train,-5])
pred2=usePL(iris[train,],iris[-train,-5])
pred3=useWhatever(iris[train,],iris[-train,-5])
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Proportions of misclassification:

```
mean(pred1==iris[-train,5])
mean(pred2==iris[-train,5])
mean(pred3==iris[-train,5])
```

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Good for choosing values for tuning parameters, or choosing between competing estimators.

R package

caret

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Also part of various techniques like Classification and Regression Tree.

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This "simulate from estimated distribution" idea is called bootstrapping.

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 $X_1,...,X_n \sim N(\mu,\sigma^2)$, with unknown μ and σ^2 .

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Take
$$\hat{\mu} = \bar{X}$$
 and $\hat{\sigma}^2 = \frac{1}{n} \sum (X_i - \bar{X})^2$.
Simulate from $N(\hat{\mu}, \hat{\sigma}^2)$.

But what is the use?

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```
orig = some available data, normal with unknown params
xbar = mean(orig)
s = sd(orig)
evt = c()
for(i in 1:10000) {
  fake = rnorm(100,mean=xbar,sd=s)
  evt[i] = abs(mean(fake)-median(fake)) > 0.05
}
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Another example: find the standard error of the 5% trimmed mean.

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But how to estimate a completely unknown distribution?

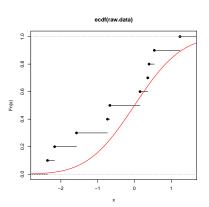
Empirical distribution

Estimate
$$F(x) = P(X \le x)$$
 by $\hat{F}_n(x) = \frac{1}{n} \# \{i : X_i \le x\}.$

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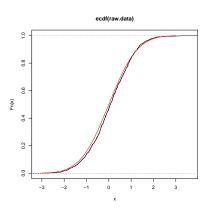
Empirical CDF (ECDF). It is a wonderful estimator. It is a step function.



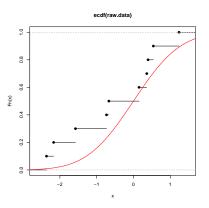
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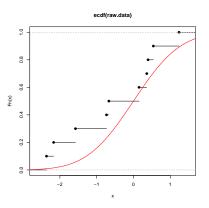


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It reduces to drawing an SRSWR from the original sample.

```
resampled.data = sample(raw.data,repl=T)
```

Bootstrap to find SE of trimmed mean

```
xbar = numeric(10000)
for(i in 1:10000) {
x = sample(raw.data,rep=T)
xbar[i] = mean(x,trim=0.1)
}
sd(xbar)
```

R package

boot

Step 4: Permutation test

This idea is mainly to be applied in testing situations.

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It does not generate data from the same process that generated the original data, but actually tries to doctor the process as well.

An example

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Consider only two types of fingerprint patterns: whorl and no-whorl.

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```
x = read.table('whorlfull.txt',head=T)
ct = table(x)
kid
mom NW W
NW 129 33
W 35 10
```

χ^2 -test

```
chisq.test(ct)
X-squared = 0.0039953, df = 1, p-value = 0.9496
The test statistic is meaningful, but the asymptotic,
null distribution needs further assumptions.
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Suppose that these assumptions are violated. Then can still perform a permutation test.

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Clearly, no association exists between mothers and children in this new data set.

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Generate lots of such data, and compute the χ^2 -statistic for each to get an idea about the null distribution.

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Generate lots of such data, and compute the χ^2 -statistic for each to get an idea about the null distribution.

Is the χ^2 -statistic value based on the actual data too large compared to these?

R code

```
mom = x[,1]
kid = x[.2]
nullchi = numeric(1000)
for(i in 1:1000) {
newkid = sample(kid)
nullchi[i] = chisq.test(table(mom,newkid))$stat
hist(nullchi,prob=T)
obschi = 0.004
mean(nullchi > obschi)
```

R package

perm, coin

Books

- Resampling Methods by Good
- Bootstrap Methods by Chernick
- ► The bootstrap, jackknife and other resampling plans by Efron