# MSA220 - Statistical Learning for Big Data Lecture 14

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Our final theme!

- When the sample size is large, there's a couple of things we need to be concerned about
- p-values become "meaningless" simply reflecting that all models are approximations of the real world
- Computations can become impossible or slow, even for simple statistical tasks

- Most methods bear a strong resemblance to stuff you're already familiar with
- Cross-validation, subsampling (bagging, RF) and bootstrap

• We will review bootstrap first because of this

An excellent book (can be found online): *An Introduction to the Bootstrap* by R. Tibshirani and B. Efron

- What and why?
- We often just trust confidence intervals that package methods spit out without thinking about underlying assumptions
- For linear models this is often OK as long as the error distribution isn't too messy or skewed and as long as the sample size is fairly big
- In other modeling scenarios, or for particular statistics of interest, we should be careful!

- In nonlinear modeling and generalized linear models, CIs provided are approximations!
- Involves linear approximations to obtain standard errors and large-sample asymptotic arguments to motivate the format of the CI

• Why not let the data do the work for us instead!

- Your observed data is a sample from the underlying population
- If you were able to repeatedly sample from this population, each time estimating your statistic of interest,  $\hat{\theta}$ ...
- then the distribution of the *theta* across the samples reflects the *sampling distribution* of this statistics

• and can be used to construct CI and for testing

#### Recap

- In the real life case you only observe one sample
- You therefore work through the properties of the statistics so that you can compute the sampling distribution statistics without having access to multiple samples
- Example: Draw x<sub>i</sub>, i = 1, · · · , n independently from population. Goal: estimate the mean μ of the population.
- $\hat{\mu} = \bar{x}$  and from iid assumption  $SE(\hat{\mu}) = \frac{1}{\sqrt{n}}\hat{\sigma}$  where

$$\hat{\sigma} = \sqrt{\sum_i (x_i - \bar{x})^2 / n - 1}$$

- If we assume  $X ilde{N}(\mu,\sigma^2)$  then  $rac{\hat{\mu}-\mu}{SE}\sim t_{n-1}$
- or even without this assumption if *n* is large we have  $\frac{\hat{\mu}-\mu}{SE}\sim N(0,1)$

• What if your statistic is more complicated so the format of the SE is unknown?

- Work it out? Use linear approximation methods (Taylor expansion, delta-method)
- Use the bootstrap!

- The idea is that you mimic the sampling from the population with a repeated sampling from the observed data
- The sampling distribution estimate you obtain by repeated sampling from the observed data can be a very good estimate of the true sampling distribution
- It doesn't always work: for extremes or "weird" statistics that are non-continuous on the true distribution or for small sample sizes.

Population



Histogram of observed sample



Histogram of bootstrap data



- For  $b = 1, \dots, B$  (B large, 1000-10000), draw a bootstrap sample from your observed data
- Non-parametric, classic: draw *n* samples with replacement.
- Alternatives: draw m < n without replacement (m-out-of-n), draw from a smooth density estimate of the data, draw from a parametric distribution
- Estimate your statistics  $\theta_b$  from each of the bootstrap data
- The distribution of θ<sub>b</sub> across b is an estimate of the sampling distribution of θ̂, the estimate from the original data

- $\bar{\theta} = \frac{1}{B} \sum_{b} \theta_{b}$  is NOT a better estimate than  $\hat{\theta}$  the purpose of bootstrap is not to improve on the estimate this way
- Bootstrap SE:  $\sqrt{\frac{\sum_{b}(\theta_{b}-\bar{\theta})^{2}}{B-1}}$  can be used to construct CI
- Bias estimate:  $\hat{\theta} \bar{\theta}$  can be used to construct a bias-corrected estimate BUT it only reflects bias with respect to estimation NOT bias induced by the wrong model assumption (that would be magic).

• Bootstrap SE:  $\sqrt{\frac{\sum_{b}(\theta_{b}-\bar{\theta})^{2}}{B-1}}$  can be used to construct CI

• Basic CI: 
$$\hat{ heta} \pm z_{1-\alpha/2}SE$$

 Note: here we are using a normal assumption for the sampling distribution BUT we could go further using the bootstrap distribution instead

- How get around the normal assumption
- Double-bootstrap
- For each bootstrap estimate  $\theta_b$ , run a second bootstrap on this bootstrap sample to obtain  $SE_b$  and compute the *pivotal* element

$$z_b = \frac{\theta_b - \hat{\theta}}{SE_b}$$

- Use the quantiles of the  $z_b$  instead of the normal quantiles
- This is called the bootstrap-t

- A conceptually simple approach is the *percentile method*
- Simple construct your confidence interval from the quantiles of the  $\theta_b$  (e.g. the 2.5% and 97.5%)!!!

• Supersimple.... BUT behaves poorly in many real-life situations.

• The rationale behind the percentile method is that for a normally distributed  $\frac{\hat{\theta}-\theta}{\sigma} \sim N(0,1)$  we have that

$$(\hat{ heta}_{lpha/2},\hat{ heta}_{1-lpha/2})$$

is a  $1-\alpha$  Cl

- With the percentile method, we assume that there is some monotone transformation g(θ̂) such that its sampling distribution is approximately normal N(0, 1)
- Why does percentile method then fail sometimes?
- There may not exist one transform that has this normalizing and variance-stabilizing effect

## BCA

- Brad Efron (The Bootstrap Guy!) proposed an improved percentile method as follows:
- Perhaps we need to correct the simple monotone transform with some bias constant and acceleration constant to make the approximate normal assumption hold

$$\phi = oldsymbol{g}( heta), \;\; rac{\hat{\phi} - \phi}{\sigma} \sim oldsymbol{N}(-z_0, 1), \; \sigma = 1 + oldsymbol{a}\phi$$

- The bias correction is obtained by the normal quantile of  $P_B(\theta_b < \hat{\theta})$
- The acceleration constant is obtained from an estimate of the skewness of the  $\theta_b$  distribution
- We adjust which quantiles in  $\theta_b$  to actually use to construct the  $1 \alpha$  Cl.
- R package boot()!!!

• New methods for dealing with large sample size

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- Parallelization or
- online updates

- This method is based on Bickel et al.'s m-out-of-n bootstrap
- m-out-of-n was shown to have better properties than regular sample-with-replacement bootstrap
- When you use m-out-of-n, you need to correct the SEs by a factor  $\sqrt{m/n}$  but otherwise it works pretty much the same way as regular bootstrap

• Here, is it used to reduce sample size!!!

- We draw s subsets of data of size m < n
- For each of the s subsets, draw r samples of size n
- Obtain point estimate and e.g. Cls from the r bootstraps

• Finally, combine the results across the s subsets

- Wait a minute! Didn't this just make the computations explode?
- Actually, no drawing a sample of size n from the subset s of size m is equivalent to assigning weights to the m observations in s
- So the computation is actually performer only on the smaller sample size *m*

### The algorithm

- For  $j = 1, \dots, s$ , draw a sample of size m (or disjoint partition of the original data)
  - For  $k = 1, \cdots, r$ ,
  - Draw weights from Multinomial(n,m)
  - Estimate your statistics of interest
- Combine by averaging quantities of interest across *s* (e.g. estimates, lower and upper CI limits, etc)

- Recommended size of  $m=n^\gamma$ ,  $\gamma\in[.5,1]$
- In the original BLB paper (Kleiner et al, 2014) they use  $\gamma = 0.6$  (reducing a data set of  $10^6$  to about 4000 for computation).
- Kleiner et al found that BLB is fairly robust to choices of *m*, consistency of estimates and good convergence rates
- Completely parallelizable for each set of size *m* so allows for fast and scalable computing

Implemented in the datadr R package

- Another variant for subsampling was proposed by Ma and Sun (2013)
- Like the BLB, they suggest that we estimate model parameters from a much smaller data set and then combine the results

• However, they differ in how the subsampling is done

• Recap from regression

• 
$$y = X\beta + \epsilon$$

• LS:  $\min_{\beta} ||y - X\beta||^2$ 

• 
$$\hat{\beta} = (X'X)^{-1}X'y$$

• 
$$\hat{y} = X\hat{\beta} = X(X'X)^{-1}X'y = Hy$$

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- Specifically,  $\hat{y}_i = \sum_j h_{ij} y_j$
- Element *h<sub>ii</sub>* is called the *leverage* of observation *i*, i.e. how much it influences its own fitted values
- Leverage basically captures if observation *i* is close or far from the center of the data. Observations near the center (in *X*-space) have limited contribution to the fit.

- Sample r observations from the original n where  $r \ll n$
- The sampling probability  $\pi_i$  for observations i is  $\pi = \frac{h_{ii}}{\sum_i h_{ii}}$
- Estimate to regression parameters
  - Alt 1: use standard OLS
  - Alt 2: use weighted LS, where the weights are the inverse sampling probabilities

- Ma and Sun found that regular OLS works better than the weighted version
- Seems simple enough!
- BUT we do need the leverage  $h_{ii}$ .
- Hm..
- The matrix *H* is *n* × *n* so we don't want to have to compute that we only care about the diagonal anyway.

• SVD to the rescue (again!)

• 
$$X = UDV'$$

• 
$$H = X(X'X)^{-1}X' = UU'$$

- and so  $h_{ii} = ||u_i||^2$  for  $u_i$  i-th row in U
- Moreover, fast randomized SVD methods exist

- Fast and simple
- A bit careful about outliers....
- A big pro: can use the subsample to visualize the data
- Model diagnostics in a big-n world and we could remove outliers at this point...

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