ECON-C4200 - Econometrics II Lecture 7: Machine learning and econometrics

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- At the end of lecture 7, you
- $1 \,$ understand what ${\bf Big} \, {\bf data}$ is
- 2 what Machine learning is
- 3 how to approach **prediction** as opposed to estimation of **(causal) parameters**
- 4 what a Ridge regression is
- 5 what a Lasso regression is
- We utilize material from Stock, J. H. & Watson, M. W. (4th Edition). *Introduction to econometrics*.

- When people talk about "Big data", they can mean many things:
 - Data with very many observations (typically, millions or more)
 - **2** Data that contain lots of variables (typically hundreds or thousands).
 - 3 Data that contain what used to be nonstandard elements such text, voice, images.

- There are numerous ways to use Big data, but by far, the most common usage is prediction in one form or the other:
 - What ads would you likely want to see (= what products might you buy)?
 - 2 How likely are you to repay your mobile phone loan? (Elisa)

3 Is this you? - recognition problems.

 Machine learning tools are also used for causal analysis both by computer scientists (the so-called Directed Acyclic Graph approach; see e.g. Paul Hunermund's MOOC) and econometricians (see e.g. Victor Chernozukov's work) but we concentrate on prediction.

What is Machine learning?

- Machine learning: using a computer and big data to "learn" (to predict as well as possible).
- Though the language used in machine learning is different from that used in econometrics, it (largely) builds on tools that are familiar to econometricians.
- A principle used (but not invented in) in machine learning is to divide data into
 - 1 a "training set" which is used to estimate the model and
 - 2 a "reserve set" (validation data / Out of sample data) which is used to compare the performance of different models.
- The objective is to predict as well as possible in the reserve set of data.
- **Caveat**: Machine learning utilizes a large variety of tools. We only cover a couple here which are directly based on regression.

- Letting go of the objective of unbiased and consistent parameters we can let go of the assumption E[ϵ|**X**] = 0.
- But we need a new assumption: Since we are using one (part of) data to estimate the model and then predict the outcomes in another (part of) data, those need to be similar (enough).
- The latter is called **O**ut-**o**f **S**ample (OOS) data (but also testing data, validation data).
- \rightarrow we assume that (X^{OOS}, Y^{OOS}) is drawn from the same distribution as the estimation data (X, Y).

- So far we have concentrated on what is needed to get unbiased and consistent estimates of β .
- In prediction, the objective is to get as accurate a prediction of the outcome as possible (in the reserve data); hence we do not care about possible biases any longer.
- However, now we need a new benchmark for what is good.
- Enter the "Oracle" who predicts as well as is possible.

• To be more practical, let us define the Mean Squared Prediction Error (MSPE) as:

$$MSPE = \mathbb{E}[Y^{OOS} - \hat{Y}(X^{OOS})]^2$$

- Y^{OOS} = outcome in the reserve / OOS-data.
- X = the variables used for prediction.
- Notice: we estimate the model using the training data, then use the predictors (X^{OOS}) in the OOS data to predict the outcome Y^{OOS}.
- Notice how MSPE is close but different from MSE.

- The **Oracle prediction** is the prediction that minimizes MSPE.
- What is this in practice?

$$Y^{Oracle} = \mathbb{E}[Y^{OOS}|X^{OOS}]$$

• Why is this? Imagine this was not the case. Then we could predict the forecast error using X^{OOS} in which case the Oracle prediction could not have been the best possible one.

• A standardized version of variable X is one that has

- mean zero and
- 2 standard deviation of one.
- It is standard in machine learning to use standardized explanatory variables.
- It is also standard to use the **demeaned** version of Y, i.e., $Y \overline{Y}$.
- \rightarrow no constant needed.
- Note: if we were interested in the coefficients, they would measure the impact on Y of a one standard deviation change in X.

Standardized regression

• A standardized regression has the same form as a regular regression:

$$Y = \boldsymbol{X}\boldsymbol{\beta} + \boldsymbol{\epsilon}$$

- This allows the use of the same "tricks" as the regular regression: polynomials, logs, interactions, ...
- **Important**: many machine learning methods (including those we cover) depend on what variables the researcher initially "proposes": e.g., they do not on their own start to introduce higher orders of a polynomial.
- By using those tricks you can increase the number of explanatory variables k to the point where k > n, i.e., you have more explanatory variables than observations.
- OLS does not work if k > n (the rank condition is not satisfied, i.e., you have more unknowns (parameters) than you have equations).

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• Let's write the true regression as

$$Y^{OOS} = \beta_1 X_1^{OOS} + \ldots + \beta_k X_k^{OOS} + \epsilon$$

• After having estimated the model with the training data we obtain $\hat{\beta}$ and can calculate

$$\hat{Y}^{OOS} = \hat{\beta}_1 X_1^{OOS} + \dots + \hat{\beta}_k X_k^{OOS}$$

- The prediction error has two sources:
 - **1** The error term e^{OOS}
 - **2** The estimation error in the parameters (coefficients): $\beta_k \hat{\beta}_k$.

- Let's denote the variance of the error term in the OOS data with $\mathbb{E}[\epsilon^{OOS}] = \sigma_{\epsilon}^2$.
- This source of prediction error will be there even if we estimate the parameters exatly, i.e., $\beta_k \hat{\beta}_k = 0$ for all k.
- $ightarrow \sigma_{\epsilon}^2$ is the MSPE of the Oracle forecast.
- What is the prediction error of our estimated model?

$$Y^{OOS} - \hat{Y}^{OOS} = (\beta_1 - \hat{\beta}_1)X_1^{OOS} + ... + (\beta_k - \hat{\beta}_k)X_k^{OOS} + \epsilon^{OOS}$$

• The MSPE of the estimated model is then

$$MSPE = \sigma_{\epsilon}^2 + [(\beta_1 - \hat{\beta_1})X_1^{OOS} + ... + (\beta_k - \hat{\beta_k})X_k^{OOS}]^2$$

• For OLS, MSPE is approximately given by

$$MSPE_{OLS} \simeq \left[1 + \frac{k}{n}\right]\sigma_{\epsilon}^2$$

- OLS has a problem: fixing the sample size *n*, MSPE is increasing in the number of predictors *k*.
- ullet ightarrow need for an estimator for which MSPE increases at a slower rate.

- It has been known for a long time (since 1950s) that by allowing bias, one can reduce *MSPE* compared to *MSPE*_{OLS}.
- With uncorrelated Xs, these estimators produce coefficients of the form

$$\hat{\beta}^{JS} = c\hat{\beta}$$

where $JS = James-Stein^1$ and 0 < c < 1.

• The Principle of Shrinkage says that we can reduce MSPE by biasing the coefficients towards zero, i.e., **shrinking** them.

¹ see James, W. & Stein, C. (1961). Estimation with quadratic loss. *Proc. Fourth Berkeley Symp. Math. Statist. Prob*, *1*, 361–379.

- One can show that in estimation, there is a trade-off between variance (of the prediction) and bias.
- This has been known for a long time, but has gained (even) more prominence with machine learning, due to its emphasis on prediction (see e.g. Giorgos Papachristoudis).
- One can show that we can rewrite the prediction error as

$$MSPE = \mathbb{E}[Y - \hat{f}(X^{OOS})]^2 = bias[\hat{f}(X^{OOS})]^2 + variance[\hat{f}(X^{OOS})] + \sigma_{\epsilon}^2$$

where $f(X^{OOS})$ is our model, e.g.,

$$f(X^{OOS}) = \beta_1 X_1^{OOS} + \dots + \beta_k X_k^{OOS} + \epsilon$$

- What happens when you increase shrinkage, i.e., decrease c?
- You increase bias by definition $\hat{\beta}^{JS} c\hat{\beta}$.
- At the same time, variance (of the prediction) decreases.
- With a large number of predictors *k*, the decrease in variance can overweigh the increase in bias.
- This would lead to a lower MSPE.
- The estimators we cover all rely on the shrinkage principle.

- 1 Split your data into estimation (training) and testing data.
- 2 Choose your model.
- **3** Estimate your model with the estimation (training) data.
- 4 Calculate the predicted values of Y, \hat{Y} , for the testing data
- Solution of the prediction errors (Y Ŷ) and dividing by n_{test}, the number of observations in the testing data.
- 6 Go back to step #2 and repeat until you cannot decrease MSPE no more.

Estimation of MSPE with cross-validation

- Best practice is to use k -fold (SW call this m-fold) cross validation.
- Example m = 10: Divide data into 10 equally sized subsamples.
- Estimate the data leaving one subsample out.
- Predict for the subsample you left out.
- Leave next subsample out, repeat.
- Repeat until you've predicted for all subsamples.
- Sum up the subsample MSPEs to get the MSPE of your estimator.

• We will cover two shrinkage - based regression approaches commonly used in machine learning:

1 Ridge regression

- 2 Least Absolute Selection and Shrinkage Operator (Lasso).
- Both **penalize** some coefficients, but do this differently.
- We use the empirical examples in SW (ch.14, newest edition).

• The principle of ridge regression is to penalize coefficients with large squared values by minimizing the following objective function:

$$S^{Ridge}(b; \lambda_{Ridge}) = \sum_{i} \left[Y - (\beta_1 X_1 + ... + \beta_k X_k)\right]^2 + \lambda_{Ridge} \sum_{j} b_j^2$$
(1)

- The second sum $\sum_j b_j^2$ is over the coefficients b_j , j = 1, ..., k.
- The term $\lambda_{Ridge} \sum_{j} b_{j}^{2}$ is the **penalty** term.
- One can show that if the regressors are uncorrelated, the ridge regression coefficients take the James-Stein form.

Ridge regression with k = 1



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The figure is for k = 1. SSR(b) = unpenalized residual sum of squares; S^{Ridge}(b) = the ridge MSPE (objective fcn.)

• While Ridge penalizes large values of squared coefficients, Lasso penalizes large coefficients with large **absolute** values:

$$S^{Lasso}(b;\lambda_{Lasso}) = \sum_{i} \left[Y - (\beta_1 X_1 + \dots + \beta_k X_k)\right]^2 + \lambda_{Lasso} \sum_{j} |b_j|$$
(2)

- The second sum $\sum_{j} |b_j|$ is over the coefficients b_j , j = 1, ..., k.
- The term $\lambda_{Lasso} \sum_{j} b_{j}^{2}$ is the **penalty** term.
- Thus Lasso and ridge very similar in appearance.



- Lasso treats large and small coefficients differently:
- It shrinks large coefficients somewhat less than ridge.
- It shrinks small coefficients to zero.
- Lasso therefore in essence "tries to get rid" of variables that affect the outcome only a little.
- A model is sparse if most of the true βs are zero and therefore

 E[Y|X] really depends on only a few of the Xs.
- Lasso seeks to produce a sparse model and therefore works well when the true model is sparse.
- This is a wonderful property when k >> n, i.e., you have many more regressors than you have observations.

Figure: LHS: large coefficients, RHS: small coefficients



- Stock and Watson also cover principal components regression.
- Briefly, the idea is to "shrink" the number of regressors using so-called **principal components** analysis.
- Once k < n, OLS can be used.

- Data: school level data on California elementary district data set with additional variables describing the schools, the students and the districts.
- There are 3 932 observations: Half of the (1 966) are used for out-of-sample prediction.
- The data has 817 predictors.

Main variables (38)

Fraction of students eligible for free or reduced-price lunch Fraction of students eligible for free lunch Fraction of English learners Teachers' average years of experience Instructional expenditures per student Median income of the local population Student-teacher ratio Number of enrolled students Fraction of English-language proficient students Ethnic diversity index	Ethnicity variables (8): fraction of students who are American Indian, Asian, Black, Filipino, Hispanic, Hawaiian, two or more, none reported Number of teachers Fraction of first-year teachers Fraction of second-year teachers Part-time ratio (number of teachers divided by teacher full-time equivalents) Per-student expenditure by category, district level (7) Per-student expenditure by type, district level (5) Per-student revenues by revenue source, district level (4)
+ Squares of main variables (38)	
+ Cubes of main variables (38)	
+ All interactions of main variables ($38 \times 37/2 = 703$)	
Total number of predictors $= k = 38 + 38 + 38 + 703 = 817$	

• Three sets of predictors are used:

- Small k = 4: Student-teacher ratio, median local income, teacher's avg. years of experience, instructional expenditures / student.
- **2** Large k = 817: See the table.
- **3** Very large k = 2065: Additional school and demographic variables, squares, cubes, interactions.
- For the Very large data set, k > n.

OLS	Ridge Regression	Lasso	Principal Components
_	-	_	_
53.6	_	-	_
52.9 1	. OLS gets wor	rse with mo	ore predictors -
Ľ			
_	2233	4527	46
78.2	39.5	39.7	39.7
64.4	38.9	39.1	39.5
_	3362	4221	69
_	39.2	39.2	39.6
$\overline{}$	39.0	39.1	39.6
	OLS 	Ridge Regression - - 53.6 - 52.9 1. OLS gets wor and you can't event - - 2233 78.2 39.5 64.4 38.9 - 3362 - 39.2 - 39.0	Ridge Regression Lasso - - - 53.6 - - 52.9 1. OLS gets worse with me and you can't even run OL - 2233 4527 78.2 39.5 39.7 64.4 38.9 39.1 - 3362 4221 - 39.2 39.2 - 39.0 39.1

Predictor Set	OLS	Ridge Regression	Lasso	Principal Components	
Small $(k = 4)$					
Estimated λ or p		2. The cross-validated MSPE.			
In-sample root MSPE	53.6	computed with the estimation			
Out-of-sample root MSPE	52.9	sample, is a good estimate of the out-of-sample MSPE			
Large (k = 817)					
Estimated λ or p	_	2233	4527	46	
In-sample root MSPE	78.2	39.5	39.7	39.7	
Out-of-sample root MSPE	64.4	38.9	39.1	39.5	
Very large (k = 2065)					
Estimated λ or p	—	3362	4221	69	
In-sample root MSPE		39.2	39.2	39.6	
Out-of-sample root MSPE	-	39.0	39.1	39.6	

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Predictor Set	OLS	Ridge Regression	Lasso	Principal Components
Small $(k = 4)$				
Estimated λ or p	-	3. Lasso, Ridge, and PC all provide big improvements over OLS		
In-sample root MSPE	53.6			
Out-of-sample root MSPE	52.9			
Large $(k = 817)$				
Estimated λ or p	_	2233	4527	46
In-sample root MSPE	78.2	39.5	39.7	39.7
Out-of-sample root MSPE	64.4	38.9	39.1	39.5
Very large $(k = 2065)$				
Estimated λ or p	_	3362	4221	69
In-sample root MSPE	-	39.2	39.2	39.6
Out-of-sample root MSPE	-	39.0	39.1	39.6

Predictor Set	OLS	Ridge Regression	Lasso	Principal Components	
Small $(k = 4)$					
Estimated λ or p	<u></u>	4. For these dat	ta, Ridge, L	asso, and PC	
In-sample root MSPE	53.6	have very similar out-of-sample MSPEs – however this will not be true in general.			
Out-of-sample root MSPE	52.9				
		For these da	ita, Ridge r	las a very sligr	π
Large $(k = 817)$		euge			
Estimated λ or p	-	2233	4527	46	
In-sample root MSPE	78.2	39.5	39.7	39.7	
Out-of-sample root MSPE	64.4	38.9	39.1	39.5	
Very large $(k = 2065)$					
Estimated λ or p		3362	4221	69	
In-sample root MSPE		39.2	39.2	39.6	
Out-of-sample root MSPE	-	39.0	39.1	39.6	

Predictor Set	OLS	Ridge Regression	Lasso	Principal Components
Small $(k = 4)$				
Estimated λ or p	_	5. For these data, there isn't much gain to using the very large data set, however this will not be true in general.		
In-sample root MSPE	53.6			
Out-of-sample root MSPE	52.9			
Large $(k = 817)$				
Estimated λ or p	_	2233	4527	46
In-sample root MSPE	78.2	39.5	39.7	39.7
Out-of-sample root MSPE	64.4	38.9	39.1	39.5
Very large $(k = 2065)$				
Estimated λ or p	_	3362	4221	69
In-sample root MSPE	_	39.2	39.2	39.6
Out-of-sample root MSPE	_	39.0	39.1	39.6



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- Many machine learning tools are based on regressions.
- They are designed for prediction, not unbiased estimation of coefficients of interest.
- Machine learning tools are especially useful when there is a large number of predictors / regressors / explanatory variables relative to the size of the data (and one only cares about prediction).
- Ridge and Lasso both utilize the shrinkage principle which builds on the bias-variance tradeoff.
- They easily outperform OLS in prediction in most cases and produce smaller MSPE.