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Artificial Intelligence in Finance

Supervised learning - continuous outcome part A

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Outline for Section 1

Introduction

Multivariate linear regression

Model description Interpretation

- Predictions
- Estimation
- Model assumptions

Model selection

In-Sample approach Selection between M-competing models Variable selection approach Out-of-Sample approach Loss functions Model confidence set

Introduction

- In supervised learning, we primarily wish to understand how a set of features is related to some future outcome(s).
- Understanding which features (and how) are related to the outcome is useful but now secondary.

Introduction

- We have at least one model that we train using a training or/and validation sample.
- Testing sample is used to evaluate (competing) models. Recall:



Training and validation (calibration) sample is used to estimate (training) and allow models to learn (validation). Most of the time they both are referred to as a training sample.

Introduction

We will learn how to go through that process:

- Key estimation models:
 - OLS.
 - LASSO, RIDGE, Elastic Net.
 - Complete subset linear regression.
 - Random forest.
 - Boosted regression trees.
- How to select and evaluate a model:
 - Standard model selection criteria (model fit, AIC, BIC, ...).
 - ML driven methods model confidence set of Hansen et al., [5].
- How to present results.

Multivariate linear regression

Outline for Section 2

Introduction

Multivariate linear regression

Model description Interpretation Predictions Estimation Model assumptions

Model selection

In-Sample approach Selection between M-competing models Variable selection approach Out-of-Sample approach Loss functions Model confidence set

Multivariate linear regression

For an overview of linear regression see Green [4]. In an **Econometrics course** the goal is to answer questions like:

- Is there a relationship between apartment area and price?
- How strong is that relationship?
- Is the relationship linear?
- How accurately can we predict price of an apartment?

In an **AI course**, we are **primarily** interested in the forecasting accuracy of one or more models. **Secondarily**, which variables are more or less important (variable importance & e**X**pl**AI**nability). Recall

Let $Y_i \in R, i = 1, 2, ..., N$, be an observed realization of an outcome variable and $X_{i,k}, k = 1, 2, ..., K$ a specific k^{th} feature:

$$E(Y|X_i) = f(X_{i,1}, X_{i,2}, ..., X_{i,K})$$

$$E(Y|X_i) = \beta_0 + \sum_{k=1}^{K} \beta_k X_{i,k}$$
 (1)

In **linear** regression, f(.) is assumed to be a **linear function**; linear in parameters β_0, β_k . Here, β_0 is referred to as an **intercept** and β_k as a **slope**.

Recall

In reality, the linear combination does not fit the data (why?):

$$Y_i \neq \beta_0 + \sum_{k=1}^{K} \beta_k X_{i,k}$$
⁽²⁾

In order to maintain equality we need to introduce a **random term** ϵ_i , with $E(\epsilon_i) = 0$, which satisfies:

$$Y_i = \beta_0 + \sum_{k=1}^{K} \beta_k X_{i,k} + \epsilon_i$$
(3)

- Why the random term ϵ_i exists?
- What does the random term represent?

Interpretation

Let P_i , i = 1, ... be the price of the i^{th} apartment and A_i apartment's area. **Interpret**:

$$P_i = 3702.5 + 3308.5A_i + \hat{\epsilon}_i \tag{4}$$



Yes, there is considerable heteroscedasticity of residuals, but point-estimates of the prediction are not affected.

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Interpretation

Let P_i , i = 1, ... be the price of the i^{th} apartment and F_i returning 1 if apartment is at the highest floor, 0 otherwise. **Interpret**:

$$P_i = 233221 + 24109.5F_i + \hat{\epsilon}_i \tag{5}$$



Interpretation

Compare more models:

- $P_i = 3703 + 3309A_i + \hat{\epsilon}_i$
- $P_i = 233221 + 24110F_i + \hat{\epsilon}_i$
- $P_i = 4202 + 3319A_i 14727F_i + \hat{\epsilon}_i$
- $P_i = -1751 + 3324A_i 15758F_i + 12545L_i + \hat{\epsilon}_i$
- $P_i = 7193 + 3197A_i + 57548F_i 960F_iA_i 31067L_i + 639L_iA_i + \hat{\epsilon}_i$

Here, L_i returns 1 if the apartment is in a building with a lift or 0 otherwise. How do you make **predictions** from such models?

Predictions

Predict the price of an apartment given the model:

 $P_i = 7193 + 3197A_i + 57548F_i - 960F_iA_i - 31067L_i + 639L_iA_i + \hat{\epsilon}_i$ (6)

The apartment has:

- area of $A_i = 50m^2$,
- is not on the highest floor $F_i = 0$,
- the building has a lift $L_i = 1$.

Predictions

The estimated model (within the training sample):

 $P_i = 7193 + 3197A_i + 57548F_i - 960F_iA_i - 31067L_i + 639L_iA_i + \hat{\epsilon}_i$ The prediction:

$$\begin{aligned} \hat{P}_i &= 7193 + 3197 \times 50 + 57548 \times 0 - 960 \times 0 \times 50 - \\ &31067 \times 1 + 639 \times 1 \times 50 \\ \hat{P}_i &= 7193 + 3197 \times 50 - 31067 + 639 \times 50 \\ \hat{P}_i &= 167926 \end{aligned}$$

Estimation

Ordinary Least **S**quares \rightarrow to estimate parameters, where $\hat{\beta}_0, \hat{\beta}_k, k = 1, 2, ...$ leads to lowest sum of squared residuals:

$$\min_{\hat{\beta}} \to \sum_{i=1}^{n} \hat{\epsilon}_{i}^{2} = \sum_{i=1}^{n} (Y_{i} - \hat{\beta}_{0} - \sum_{k=1} \hat{\beta}_{k} X_{i,k})^{2}$$

How do we find the minimum of this function?

- There are **other possible criteria** not just squared errors.
 - Absolute least squares.
 - Weighted least squares with various weighting schemes.

$$\min_{\hat{\beta}} \to \sum_{i=1}^{n} \hat{\epsilon_i}^2 = \sum_{i=1}^{n} w_i (Y_i - \hat{\beta}_0 - \sum_{k=1}^{n} \hat{\beta}_k X_{i,k})^2$$

■ You can be **quite creative** about weights you employ.

Estimation

We estimate β 's, but also the standard error (variation of) β , $var(\beta)$:

- It allows us to assess the **accuracy** of the parameter estimate.
- It allows us to test (under certain assumptions) statistical hypotheses about β, e.g. H₀ : β₁ = 0.
- We can assess what influences the precision of our parameter estimates.
- Given initial assumptions about the model (see later) we have many approaches to estimate var(β).
- Opposed to econometrics, in machine learning, we do not care so much about the variance of parameter estimates, instead we care about the variance and bias of our predictions. Although the two are related.

Model assumptions

Model should work best (including inferences) after satisfying a series of assumptions/rules (see Greene [4]):

- 1. Model is **linear in parameters** β .
- 2. Independent variables are not stochastic.
- 3. $E(\epsilon_i|X_i) = 0 \rightarrow E(Y_i|X_i) = \beta_0 + \sum_{k=1} \beta_k X_{i,k}$.
- 4. Homoscedasticity of residuals $var(\epsilon_i|X_i) = \sigma^2$.
- 5. For $i \neq j$, residuals ϵ_i and ϵ_j are **uncorrelated** (time-series, spatial,...).
- 6. There is no covariance between ϵ_i and X_i .

Model assumptions

Model should work best (including inferences) after satisfying a series of assumptions/rules (see Greene [4]):

- 7. Number of observations needs to be at least the number of parameters, $K \le N$.
- 8. Variance of X_k should be finite and positive.
- 9. Regression model should be well-specified.

Assumption No. 9 is quite important and the reason, why artificial intelligence methods (statistical learning in general) will likely get the upper-hand in the future.

Model selection

Outline for Section 3

Introduction

Multivariate linear regression

Model description Interpretation Predictions Estimation Model assumptions

Model selection

In-Sample approach

Selection between M-competing models Variable selection approach

Out-of-Sample approach

Loss functions Model confidence set

Model selection

'All models are wrong, but some are useful.' Box, G. E. P. (1979, [3])

Standard econometric assumption suggests that we have a well-specified regression.

- That is **almost surely not true**.
- In Finance (observational studies), we can be pretty sure we do not have a well-specified model.

Instead of assuming that one model represents the 'correct specification', an **AI (data)-driven approach** wants to **learn** which model(s) tend to perform better.

In-Sample approach

In an in-sample approach, we do not learn from new data, but select a model (or a variable) using the **training dataset** (not validation) **only**. We will consider following scenarios:

- Select between M-competing models.
 - **R**² and adjusted R^2 .
 - Model confidence set of Hansen et al., (2011, [5]).
 - Akaike and Schwartz Information Criteria.
- Variable selection approach.
 - Backward selection.
 - Forward selection.
 - Step-wise selection.

Example:

Let $R(1)_i, R(2)_i, ..., R(5)_i$ denote the number of rooms of i^{th} apartment and FR_i, PR_i, NW_i a full- and partial-reconstruction and a new apartment. We have M = 4 competing models:

$$P_{i} = \beta_{0} + \beta_{1}A_{i} + \epsilon_{i}$$

$$P_{i} = \beta_{0} + \beta_{1}A_{i} + \sum_{k=2} \beta_{k}R(k)_{i} + \epsilon_{i}$$

$$P_{i} = \beta_{0} + \beta_{1}A_{i} + \beta_{2}FR_{i} + \beta_{3}PR_{i} + \beta_{4}NW_{i} + \epsilon_{i}$$

$$P_{i} = \beta_{0} + \beta_{1}A_{i} + \beta_{2}FR_{i} + \beta_{3}PR_{i} + \beta_{4}NW_{i} + \sum_{k=5} \beta_{k}R(k)_{i} + \epsilon_{i}$$

Simple procedure for each of the measures:

- 1. Estimate all *M* models using data from the training sample.
- 2. For each model, **calculate** adjusted R^2 (AIC and/or BIC).
- 3. **Select** the model that has highest R^2 (lowest AIC and/or BIC).

Model fit - R²

Let Y_i , i = 1, 2, ..., N be the outcome variable and let *TSS* denote the Total Sum of Squares, i.e. the variability we want to explain in the first place:

$$TSS = \sum_{i=1}^{N} (Y_i - \bar{Y})^2$$
 (7)

We can decompose this variability into the part that we can explain via our model, the Explained Sum of Squares and into the part that the model was not able to explain, the Residual Sum of Squares:

$$TSS = ESS + RSS = \sum_{i=1}^{N} (Y_i - \bar{Y})^2 = \sum_{i=1}^{N} (\hat{Y}_i - \bar{Y})^2 + \sum_{i=1}^{N} (Y_i - \hat{Y}_i)^2$$
(8)

Model fit - R²

Visual representation of the **components of** the R^2 :



$$TSS = ESS + RSS = \sum_{i=1}^{N} (Y_i - \bar{Y})^2 = \sum_{i=1}^{N} (\hat{Y}_i - \bar{Y})^2 + \sum_{i=1}^{N} (Y_i - \hat{Y}_i)^2$$
(9)

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Model fit - R²

Recall:

$$TSS = ESS + RSS$$

TSS is given by the outcome variable (Y_i). We can influence ESS (RSS) by using a suitable model. The lower the RSS, the better the model, which leads us to the following, **coefficient of determination**:

$$0 \leq R^2 = 1 - \frac{RSS}{TSS} \leq 1 \tag{10}$$

 R^2 measures the ratio of explained variability of the outcome variable. It is one of the most popular measures of a goodnes of a model; in general, not just for linear regressions.

- With the increase of features in the linear regression model, the R^2 almost surely increases.
- It is still very informative as it tell us how one model (feature) increases our understanding about the variation of the outcome variable.
- We know that having too much parameters leads to over-fitting → we should penalize more complex models; principle of parsimony.

Let *p* denote the number of features, the adjusted R^2 is given as:

$$R_a^2 = 1 - \left(\frac{n-1}{n-p}\right) \left(1 - R^2\right) \le 1 \tag{11}$$

Let LL(m) be the log-likelihood (see Econometrics course) of a model \rightarrow how good the model fits the data. The Akaike (1974, section V in [1]) and Schwartz (1978, [6]) information criteria are given as:

$$AIC = -2LL(m) + 2p$$

$$BIC = -2LL(m) + ln(N)p$$
(12)

Specifically if comparing models from linear regressions, assuming $\epsilon_i \sim N(\mu, \sigma^2)$, the term -2LL(m) is replaced with:

$$Nln\left(\frac{RSS_m}{N}\right) + c$$
 (13)

and constant *c* can be ignored.

Let's turn back to our example:

Comparing models						
Criteria	OLS-1	OLS-2	OLS-3	OLS-4		
<i>R</i> ²	61.92%	63.51%	64.89%	66.13%		
R_a^2	61.91%	63.45%	64.85%	66.05%		
AIC	89154	89016	88879	88763		
BIC	89173	89065	88915	88831		

OLS - 4 seems to be the preferred according to all three criteria.

Backward elimination:

- 1. Select a threshold p-value (say 0.10).
- 2. Estimate regression with all features.
- 3. Estimate p-value (H_0 : $\beta = 0$, and use appropriate technique to estimate coefficient standard errors, e.g. bootstrapping, HC, HAC).
- 4. Find the feature with the highest p-value and if it is above the threshold, remove the feature from the specification.
- 5. Estimate the model with new specification.
- 6. Repeat step 3 to 5 until the highest p-value is less than the threshold p-value.

Variable (and model) selection approach

Backward elimination: Looking back at our example, the selected model (OLS-B henceforth) is:

$$P_i = \beta_0 + \beta_1 A_i + \beta_2 R(2)_i + \beta_3 R(3)_i + \beta_4 R(5)_i + \beta_5 F R_i + \beta_6 N W_i + \epsilon_i$$

with parameter estimates given by:

$$P_i = 1266 + 3123A_i - 10016R(2)_i - 30866R(3)_i + 72638R(5)_i + 31793FR_i + 63849NW_i + \epsilon_i$$

Compared to m = 4, the $R_a^2 = 66.06\%$, *AIC* = 88758 and *BIC* = 88808 are now improved.

Forward elimination:

- 1. Select a threshold p-value (say 0.10).
- 2. Estimate a regression with an intercept.
- 3. Add one feature which leads to the lowest p-value on the corresponding regression coefficient ($H_0: \beta = 0$).
- Repeat step 3 until you are **unable to find** a feature that would have a *p*-value below the threshold defined in step 1.

Forward elimination: Looking back at our example, the selected model (OLS-F henceforth) is:

$$P_i = \beta_0 + \beta_1 A_i + \beta_2 R(1)_i + \beta_3 R(2)_i + \beta_4 R(3)_i + \beta_5 SR_i + \beta_6 NW_i + \beta_7 FR_i + \beta_8 R(5)_i + \epsilon_i$$

with parameter estimates given by:

$$P_i = 5094 + 3094A_i - 7282R(1)_i - 13536R(2)_i - 33735R(3)_i + 1856SR_i + 65418NW_i + 33055FR_i + 72705R(5)_i + \epsilon_i$$

The $R_a^2 = 66.05\%$, *AIC* = 88761 and *BIC* = 88823 are **worse** as for m = 5 (OLS-B).

Bi-directional selection

- 1. Select a threshold *p*-value (say 0.10).
- 2. Estimate a regression with an intercept.
- 3. Add one feature to the model that leads to the lowest *p*-value.
- 4. Remove feature(s) that have a *p*-value above the threshold.
- 5. Repeat step 3 to 4 until adding a new feature does not lead to feature's *p*-value below the threshold *p*-value.

Some variables may be added at one iteration, removed in another and added back to the set of preferred features latter on!

Bi-directional elimination: Looking back at our example, the selected model is:

$$P_i = \beta_0 + \beta_1 A_i + \beta_2 R(2)_i + \beta_3 R(3)_i + \beta_4 NW_i + \beta_5 FR_i + \beta_6 R(5)_i + \epsilon_i$$

with parameter estimates given by:

$$P_i = 1266 + 3123A_i - 10016R(2)_i - 30866R(3)_i + 63849NW_i + 31793FR_i + 72638R(5)_i + \epsilon_i$$

The model is the same as m = 5!

Forecasts errors lead to costs:

- **Economic losses** are very specific and depend on the application of the forecast.
- Statistical losses are more general. For continuous variables most common are:
 - Square Error.
 - Absolute Error.

We use these functions to **optimize** (parameters) our models and to evaluate model forecasts - to **rank** models.

Let Y_i , i = 1, 2, ...N, be the i^{th} observed outcome variable and $\hat{Y}_{i,m}$, m = 1, 2, ... the corresponding forecast from model m. The two common loss (cost) functions are:

Mean Square Error.

$$MSE_m = N^{-1} \sum_{i=1}^{N} (Y_i - \hat{Y}_{i,m})^2$$
(14)

Mean Absolute Error.

$$MAE_m = N^{-1} \sum_{i=1}^{N} |Y_i - \hat{Y}_{i,m}|$$
 (15)



- Different preference for larger losses.
- Symmetric behavior for under/over predictions.
- You can specify your own losses (e.g. asymmetric losses).

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Let's continue with out example. Using models from the **training** sample only (not validation) we predict prices in the **testing** sample: **training** \rightarrow **testing**.

Comparing models					
Models	MSE ×10 ⁻⁹	MAE $\times 10^{-4}$			
OLS-1	12.31	6.07			
OLS-2	11.99	5.93			
OLS-3	11.37	5.64			
OLS-4	11.16	5.56			
OLS-B	11.18	5.56			
OLS-F	11.17	5.56			

In the previous example, differences between models m = 4, 5, 6 seem to be negligible. Perhaps the three models lead to similarly accurate forecast.

Can we perform a statistical test to compare models?

Currently popular is the **Model Confidence Set** (MCS henceforth) approach of Hansen et al., (2011, [5]). We will follow the exposition of Bernardi and Catania (2018, [2]).

Let \hat{M}^0 denote an initial set of q models (m = 1, 2, ..., q). The goal is for a given confidence level of $1 - \alpha$ arrive at a smaller set of superior models denoted as $\hat{M}^*_{1-\alpha}$ with $q^* \leq q$. The general algorithm [2]:

- 1. Set $M = M_0$, i.e. the initial set consists of all models.
- 2. Test a null hypothesis that all models lead to equal predictive accuracy (EPA).
 - If the hypothesis is not rejected, the algorithm is terminated.
 - If the hypothesis is rejected, determine the worst model and remove it from *M*.
- 3. Repeat step 2 until the EPA hypothesis is not rejected.

The **EPA hypothesis**. Let $l_{m,i}$ be a loss (e.g. squared error) of the m^{th} model m = 1, 2, ..., q at predicted observation *i*. The **loss differential** is:

$$d_{m,r,i} = l_{m,i} - l_{r,i}; \ m \neq r$$
 (16)

The EPA hypothesis for a given set of models *M* can be defined as:

$$\begin{aligned} & H_{0,M}: E(d_{m,r}) = 0, \quad \forall m, r = 1, 2, ..., q \\ & H_{1,M}: E(d_{m,r}) \neq 0, \quad \text{for some } m, r = 1, 2, ..., q \end{aligned}$$

For comparing only two models, the **test statistics** is given by:

$$t_{m,r} = \frac{\bar{d}_{m,r}}{\sqrt{\hat{var}(\bar{d}_{m,r})}}$$
(18)

where $\overline{d}_{m,r}$ is the average loss between models *m* and *r*. The challenge is to calculate $\widehat{var}(\overline{d}_{m,r})$. Hansen et al (2011, [5]) recommends block bootstrap procedure (can be adapted for cross-sections). Comparing multiple models, the **test statistics** becomes:

$$T_{RM} = max_{m,r\in M}|t_{m,r}| \tag{19}$$

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Let's continue with out example. Using models from the **training** sample only (not validation) we predict prices in the **testing** sample: **training** \rightarrow **testing**.

Comparing models						
Models	$MSE imes 10^{-9}$		$MAE imes 10^{-4}$			
OLS - 1	12.31		6.07			
<i>OLS</i> – 2	11.99		5.93			
OLS — 3	11.37	†	5.64			
<i>OLS</i> – 4	11.16	†	5.56	†		
OLS — B	11.18	†	5.56	†		
OLS - F	11.17	†	5.56	†		

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Model selection Out-of-Sample approach

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Supervised learning - continuous outcome part A

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