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MF793 - Fall 2021

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OLS: When things go wrong

- A. Quick Return on Correlated Regressors
- B. Omitted X variables
- C. RHS variable measured with error
- D. Using the residuals as diagnostic of the (unknown) noise properties
- E. Difference between residual and forecast error, the leave-one-out residual
- F. Errors are not i.i.d: Heteroskedastic and / or correlated errors,

Readings: Hansen Ch. 2, 3, 4

Greene: Ch. 4.7

A. Quick return on correlated regressors

Recall
$$V(\hat{\beta}_{OLS}) = \sigma^2 (X'X)^{-1}$$

Highly correlated X variables increase the variance of the OLS. .. Which coefficient?

Is there some intuition in (X'X)⁻¹ diagonal elements?

Yes, we can show (no proof) that:

$$V(\widehat{\beta}_{\text{OLS},k}) \equiv \sigma^2 \{ (X'X)^{-1} \}_{kk} = \frac{\sigma^2}{\left(1 - R_k^2\right) \sum_{1}^{T} (x_{k,i} - \overline{x_k})^2}$$

Where:

 \mathbb{R}^2_k is the R2 of the regression of x_k on all the other x's

 $\frac{1}{(1-R_k^2)}$ is called the **variance inflation factor** of the estimate of β_k .

B. Missing Variable, we regress Y on X₁, forget a variable X₂

B1 Effect on the coefficient estimate

Say the correct model is:

$$Y = X_1\beta_1 + X_2\beta_2 + \varepsilon^*, E(\varepsilon^* \mid X_1, X_2) = 0$$
 [1]

But we think **incorrectly** that:

$$Y = X_1 \beta + \varepsilon$$
,

[2]

Our estimate from the wrong model is: $b = (X_1'X_1)^{-1} X_1'Y$

$$b = (X_1'X_1)^{-1} X_1'Y_1$$

When is b an unbiased estimate of β_1 ?

$$b = (X_1'X_1)^{-1} X_1' \mathbf{Y} = (X_1'X_1)^{-1} X_1' (X_1\beta_1 + X_2 \beta_2 + \varepsilon)$$
 <- Y by true model

$$= \beta_1 + (X_1'X_1)^{-1} X_1' X_2 \beta_2 + (X_1'X_1)^{-1} X_1' \varepsilon$$

$$= \beta_1 + (X_1'X_1)^{-1} X_1' \varepsilon + (X_1'X_1)^{-1} X_1' X_2 \beta_2$$

$$E(\mathbf{b}) = \beta_1 + (X_1'X_1)^{-1} X_1' X_2 \beta_2$$

 $(X_1'X_1)^{-1} X_1'X_2$: The matrix of the regression coefficients X_2 on X_1

The estimate b is biased iff X_1 is uncorrelated with X_2 => That is if the regression coefficient of X_2 on X_1 is zero.

B2 Risk of too many vs Risk of not enough variables

- Omitting variables:
 - \circ $\hat{\beta}$ for the included variables can be a biased estimate of β_1 We can also show that variance of the estimate is smaller (than if using correct model)! no proof
 - \circ Effect of omitting X2 on a forecast \hat{y} is less obvious
 Biased coefficient estimate in effect compensates (not totally) for missing variable
- Do we run the regression to test a hypothesis on β_1 or to forecast y?
- Contrast with including too many "not so useful" variables! If the included variables are (highly) correlated
 - o variance of the coefficient estimator increases
 - o variance of the forecast error increase
 - o No bias
 - o The inference is correct, both for estimation and forecast
- Boils down to modeling strategy: risk of too many variables vs risk of missing essential variables.
 - o Old days: Bottom up modeling favored, start with few X variables, build up the model
 - o Less old days: Top down preferred, start from many variables, reduce the model.
 - o Either way is deeply flawed.

Both introduce severe biases known as "pretest biases" or "data mining biases"

- What are data mining biases:
 - Say your final model is: Y on X₁
 - \circ Your final model won after many trials of Y on X_1 .. , and on X_2 , X_3 , .. which were eliminated.
 - o Standard t-test for $\hat{\beta}_1$ in your final model **overestimates** its significance.

Think: Assume you start with 100 candidate X variables, trying to forecast Y

- What to do?
 - o No mechanical variable selection technique is reliable ...
 - ... without a solid dose of common sense and subject matter knowledge.
 - Know your subject, don't use mechanical model selection techniques
 - If you must do model selection, avoid criteria that lead to large models.
 R² is useless for model selection, adjusted R² is hardly better.
 Prefer AIC or SIC criteria (will see in Time series)
 - If you did any model selection, you can not use the usual standard error and t-statistic for your estimates and forecasts.
 - because the data-mining you conducted modified the "nominal" distribution of your test
 - Even better: model combination rather than model selection, odds ratios.
 - Why "throw away" models?
 - Does a portfolio manager throw away the stocks with higher variance?

C Variables are measured with errors

C1 Dependent variable Y measured with error

• Theoretical Model: $Y^* = X\beta + \varepsilon$

We don't know Y*, we **measure** it: $Y = Y^* + v$ v: random measurement error

So the model for the data Y is: $Y = X\beta + \varepsilon + v$ [1]

• If the measurement error is additive, unrelated to X, and homoskedastic,

the only consequence is a higher variance of noise when we use Y

=> lower fit.

But the model's assumption are maintained, the inference is correct.

C2 Independent variable X measured with error

• Theoretical Model: $Y = \alpha + X^*\beta + \epsilon$

We don't know X*, we measure it: $X = X^* + v$ $v \sim (0, \sigma_v)$

So the model for the data X,Y is: $Y = \alpha + X\beta + (\epsilon - \beta \nu)$ [2]

• Cov(ε - βv , X) = Cov(ε , X) - β Cov(ν , X) = 0 - β Cov(X*+ ν , ν) = - β Var(ν) \neq 0

- Even for the simplest assumption on the measurement error: $Cov(v, X^*) = Cov(v, \varepsilon) = 0$, the errors in the model we run ([2]) are correlated with the RHS variable we use, X. The OLS estimate is biased.
- We can also show that the bias of the OLS estimate of β

no proof

- o is generally toward zero,
- $\circ~$ is larger the larger the size of measurement error σ_{ν} .

When the RHS variables are estimated with error, the OLS beta estimator is generally biased toward zero, the bias is larger the larger σ_{ν}

D Using the residuals to conduct model diagnostic

- We use the residuals *e* to diagnose
- 1) patterns of dependencies between X and ϵ
- 2) heteroskedasticity of the noise
- 3) patterns of dependencies among ϵ
- Let's understand the distribution of these residuals
- Covariance matrix of the residuals:

Recall
$$e = M\epsilon$$
 where $M = I - X(X'X)^{-1}X' = I - P$ M is symmetric $V(e) = E(ee') = E(M \epsilon \epsilon' M') = M E(\epsilon \epsilon') M = \sigma^2 M \neq \sigma^2 I$ [1]

When the true noise is i.i.d., the residuals

are **not uncorrelated** with one another and **not homoskedastic!**

Variance of one residual:

$$V(e_i) = \sigma^2 m_{ii} = \sigma^2 (1 - p_{ii})$$
 [1']

• Result: Standardized residuals can not be computed as $\frac{e_i}{s}$ even if s^2 is an unbiased estimate of σ^2 .

• Standardized residuals must be computed as

$$\frac{e_i}{s\sqrt{1-p_{ii}}}$$
 [2]

Residuals as in [2] are sometimes called **studentized** residuals.

Be careful when reading / using different packages!

So, if we call [2] a standardized residual, what is a studentized residuals?

It is a (third!) standardization where s in [2] is replaced by s_i , an estimate of σ which excludes the residual i. ... That is we leave observation i out to compute s_i .

There is no common standard used by everybody to refer to the standardized (or studentized?) residuals!

• Never mind the jargon, you should know [1'] and use [2]. How is it done in R:

influence(model)\$hat

gives the vector of diagonals values of P, the Hat matrix

plot(model)

gives diagnostic plots based on standardized residuals computed with [2]

• It's a good time to discuss "leave one out" diagnostics.

E. Leave-one-out and all that sort of things. (Hansen 3.16 for details)

• What is the difference between a residual and a forecast error?

Residual: $e_i = y_i - x_i \hat{\beta} = y_i - \hat{y}_i$ Forecasting error: $e_f = y_f - x_f \hat{\beta} = y_f - \hat{y}_f$

Observation *i* was used to estimate β

Observation f was not!

- Say we use T observations to estimate β , and we have T₂ additional "out-of-sample" observations.
 - \circ We can compute an **Out-Of-Sample R**² with the T₂ observations which **did not** influence $\hat{\beta}$
 - o OOS R² much lower than the in-sample R² indicates model instability ... or data mining.

Save data for OOS forecasting, as diagnostic of stability / data mining on your model.

- The **Leave-one-out** technique.
 - ο For each (x'_i, y_i) , estimate β by **excluding** observation I, $-> \widehat{\beta}_{-\iota}$, \widetilde{y}_{ι} , \widetilde{e}_{ι} **Leave-one-out**
 - o Can show: $\hat{y}_i \tilde{y}_i = p_{ii} \tilde{e}_i$. Diagonal elements of the P matrix diagnose influence point.
 - o Compute R² with the leave-one-out fitted values and residuals.
 - \circ The **leave-one-out** \mathbb{R}^2 is a better model selection criterion than \mathbb{R}^2 and \mathbb{R}^2
- Cross-validation

Large models vastly **over**fit: They achieve very high in-sample R² which collapse out of sample.

Say: T observations, Leave K out, estimate on T-K

Only 1 way to leave-one-out, MANY way to leave k out

Cross validation is a generalization of "leave one out"

F Heteroskedasticity, Correlated errors, GLS or adjusted OLS Standard errors?

F1. Complication:
$$E(\epsilon\epsilon)$$

$$E(\varepsilon \varepsilon') \neq \sigma^2 I$$
 $E(\varepsilon \varepsilon') = \sigma^2 \Omega = \Sigma$

- We write $\sigma^2 \Omega$ so that the general model nests the iid model. But of course in many situations, σ is not identified uniquely. It's only a convention.
- Heteroskedasticity: $E(\varepsilon \varepsilon') = D \neq \sigma^2 I$, $E(\varepsilon_i^2) \neq E(\varepsilon_i^2)$

$$E(\varepsilon\varepsilon') = D \neq \sigma^2 I$$

$$E(\varepsilon_i^2) \neq E(\varepsilon_j^2)$$

$$\sigma^2 \mathbf{\Omega} = \sigma^2 \begin{bmatrix} \omega_1 & 0 & \cdots & 0 \\ 0 & \omega_2 & \cdots & 0 \\ & & \vdots \\ 0 & 0 & \cdots & \omega_n \end{bmatrix} = \begin{bmatrix} \sigma_1^2 & 0 & \cdots & 0 \\ 0 & \sigma_2^2 & \cdots & 0 \\ & & \vdots \\ 0 & 0 & \cdots & \sigma_n^2 \end{bmatrix}.$$

Examples:

- Error variance proportional to one of the regressors: \(\sigma_i = \sigma \sqrt{x_i} \)
 Error variance ... varies with time or previous shock: \(\sigma_t = \alpha + \delta r_{t-1}^2 \) ARCH model

• Correlated errors $E(\varepsilon_{\rm I} \, \varepsilon_{\rm j}) \neq 0$ for $i \neq j$, $E(\varepsilon \varepsilon') = \sigma^2 \, \Omega$

For homoskedastic, correlated errors, σ^2 is identified. It the constant error variance and the Ω matrix is a correlation matrix.

• Examples:

o Cross-sectional regression where observations indicate for example industry / region

Group effects: $i \neq j$ in same industry/region, $E(\varepsilon_i \varepsilon_j) \neq 0$

 $i\neq j$ in different industries/regions $E(\epsilon_i\epsilon_j) = 0$

o Time Series regression

Autocorrelation of order 1: $\varepsilon_t = \rho \ \varepsilon_{t-1} + \nu_t$ AR(1)

$$\sigma^{2}\mathbf{\Omega} = \sigma^{2} \begin{bmatrix} 1 & \rho_{1} & \cdots & \rho_{n-1} \\ \rho_{1} & 1 & \cdots & \rho_{n-2} \\ & \vdots & & \vdots \\ \rho_{n-1} & \rho_{n-2} & \cdots & 1 \end{bmatrix}$$

where ρ_{K} = Corr(ϵ_{t} , ϵ_{t-k}). We expect ρ to be higher if k is smaller (close observations) We will see that for the AR(1) model $\rho_{K} = \rho^{K}$

F2. General Solution: Generalized Least Squares (GLS)

- Matrix Result: $\forall \Omega, \exists P / \Omega = P' D P$ [1] P'P = P P' = I and D is (the) diagonal (matrix of Eigen values).
- Pre-multiply Ω by $P^* = D^{-0.5}P$

$$D^{-0.5}$$
 P Ω P' $D^{-0.5}$ = $D^{-0.5}$ P P'DP P' $D^{-0.5}$ = $D^{-0.5}$ D $D^{-0.5}$ = I_T

P* Ω P*' = I_T Keep this ready near the stove.

- Pre-multiply our regression: $P^* Y = P^* X \beta + P^* \epsilon$ $E(\epsilon \epsilon') = \sigma^2 \Omega$ [2]
- $E(\varepsilon^*\varepsilon^{*'}) = E(P^*\varepsilon\varepsilon'P^{*'}) = P^*\sigma^2\Omega$ $P^{*'} = \sigma^2I$ => OLS applies to [2], it is BLUE for [2]
- GLS is just OLS applied to the transformed system [2]

$$\hat{\beta}_{GLS} = (X^{*'} X^{*})^{-1} X^{*'} Y^{*} = (X'\Omega^{-1}X)^{-1} X'\Omega^{-1} Y$$

$$V(\hat{\beta}_{GLS}) = \sigma^{2} (X^{*'}X^{*})^{-1} = \sigma^{2} (X'\Omega^{-1}X)^{-1}$$
[4]

- Are we done? No! Not even started! We need an estimate of Ω to plug into [3][4]: $\widehat{\Omega}$ to get a *Feasible GLS*
- Ω is a TxT matrix.

Under i.i.d. errors, we reduced it to one parameter, $\sigma^2 I_T$.

We now have T(T+1)/2 parameters and only T observations

• Can't estimate T(T+1)/2 parameters with T observations!

Need to specify a model of correlation / heteroskedasticity to reduce the number of parameters

Examples: autocorrelated residuals with AR(1) One parameter ρ describes entire matrix

Weighted Least Squares, $\sigma_i = \sigma x_i$ Observable describes variance

Noise variance follows a GARCH process

• **Problem:** GLS is very sensitive to assumptions on the covariance matrix

Wrong $\widehat{\Omega}$ means wrong \widehat{eta}_{GLS} : GLS can then be biased and inefficient

The GLS variance covariance matrix is then incorrect: $V(\hat{\beta}_{GLS}) \neq \sigma^2 (X'\Omega^{-1}X)^{-1}$

GLS is a nice idea but you need to be confident that your model is well specified.

F3. Remain with OLS, get a good estimate of the OLS standard error. HAC standard errors

What is wrong with OLS if $E(\varepsilon \varepsilon') = \sigma^2 \Omega$?

- OLS is likely inefficient.
- But we are not sure of what model to use for GLS.
- The variance of $\hat{\beta}_{OLS}$ is **not** $\sigma^2(X'X)^{-1}$.

That formula is wrong, bye-bye confidence intervals for estimates and forecasts

Possible view:

- GLS is tricky: if we pick the wrong Ω , the remedy may be worse than the disease.
- Better a somewhat inefficient estimator (OLS) with properly computed standard errors than a mis-specified estimator with incorrectly computed everything.

Counter View:

At least for Time-Series models, one can easily devise a GLS based AR model + GARCH model for the errors. It would likely dominate OLS.

This is true: We will do this in MF840

• For now, we introduce the **robust standard errors** for:

Unspecified heteroskedasticity Hal White (1983) Unspecified autocorrelation Newey West (1987).

They allow us to keep using OLS while correcting the variance formula

• For heteroskedasticity: $E(\epsilon \epsilon') = D$. We don't know D!

$$V(\hat{\beta}_{OLS}) = E[(X'X)^{-1} X' \varepsilon \varepsilon' X (X'X)^{-1}] = \sigma^2 (X'X)^{-1} X' D X (X'X)^{-1}$$

Ideal estimator if we knew the true noises: $X' \varepsilon \varepsilon' X = \sum_i x_i x_i' \varepsilon_i^2$

Hal White (1983):

We can estimate $E(X' \in \mathcal{E}' X)$ consistently even if we can't estimate D consistently

We replace the true ε_i 's with the residual e_i 's or better, rescaled e_i 's.

Of course, it does not estimate D consistently, $\{diag(e^{2}_{i})\}$, one observation per parameter!

But it estimates X' D X and hence $V(\hat{\beta}_{OLS})$ which are KxK

• Autocorrelated Errors: Hansen-Hodrick and Newey and West, propose similar strategies. They use ad-hoc estimates of autocorrelation until fairly high lags

Hansen-Hodrick worked in a context of overlapping observations where they **knew** how many lags of autocorrelations were in the data.

In general, it can create non-positive covariance matrices. Newey-West corrects this problem and is preferred.

- These are called **HAC standard errors**: Heteroskedasticity Autocorrelation Consistent OLS may not be the "best" estimator but you can specify its uncertainty properly
- In R: command coeftest, needs packages sandwich and lmtest.