Foundations of Linear Regression

2. Review, Properties and Assumptions, Matrix Form

GOVT 6029 - Spring 2019

Cornell University

Outline

1. Housekeeping

2. Main ideas

- 1. Correlation coefficient describes the strength and direction of the linear association between two numerical variables
 - 2. Least squares line minimizes squared residuals
 - 3. Interpreting the least squares line
 - 4. Predict, but don't extrapolate
 - Regression in social sciences
 - Distributions
 - 7. Linear Regression in Matrix Form

3. Assumptions

Announcements

• Problem Set 1 ... Questions, comments?

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- Due Friday html and Rmd

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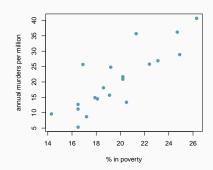
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Your turn

Which of the following is the best guess for the correlation between annual murders per million and percentage living in poverty?

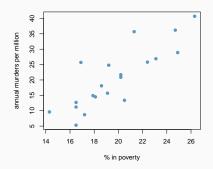
- (a) -1.52
- (b) -0.63
- (c) -0.12
- (d) 0.02
- (e) 0.84



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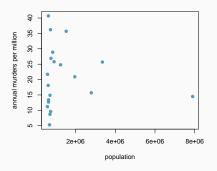
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Your turn

Which of the following is the best guess for the correlation between annual murders per million and population size?

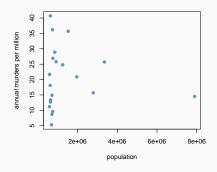
- (a) -0.97
- (b) -0.61
- (c) -0.06
- (d) 0.55
- (e) 0.97



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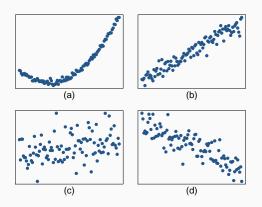
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Assessing the correlation

Your turn

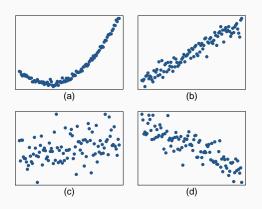
Which of the following is has the strongest correlation, i.e. correlation coefficient closest to +1 or -1?



Assessing the correlation

Your turn

Which of the following is has the strongest correlation, i.e. correlation coefficient closest to +1 or -1?



 $(b) \rightarrow$ correlation
means <u>linear</u>
association

Play the game!

http://guessthecorrelation.com/

Spurious correlations

Remember: correlation does not always imply causation!

http://www.tylervigen.com/

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(2) Least squares line minimizes squared residuals

 Residuals are the leftovers from the model fit, and calculated as the difference between the observed and predicted y:

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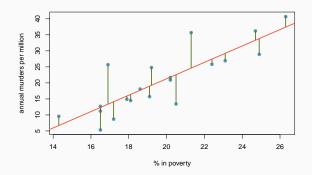
- The least squares line minimizes squared residuals:
 - Population data: $\hat{y} = \beta_0 + \beta_1 x$
 - Sample data: $\hat{y} = b_0 + b_1 x$

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Interpreting the last squares line

• *Slope:* For each <u>unit</u> increase in <u>x</u>, <u>y</u> is expected to behigher/lower on average by the slope.

$$b_1 = \frac{s_y}{s_x} R$$

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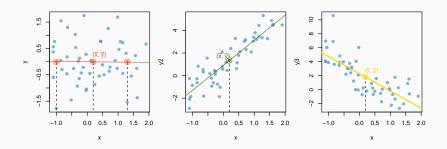
$$b_0 = \bar{y} - b_1 \bar{x}$$

• The calculation of the intercept uses the fact the a regression line **always** passes through (\bar{x}, \bar{y}) .

• If there is no relationship between x and y ($b_1 = 0$), the best guess for \hat{y} for any value of x is \bar{y} .

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Your turn

What is the interpretation of the slope?

$$\widehat{murders} = -29.91 + 2.56$$
 poverty

- (a) Each additional percentage in those living in poverty increases number of annual murders per million by 2.56.
- (b) For each percentage increase in those living in poverty, the number of annual murders per million is expected to be higher by 2.56 on average.
- (c) For each percentage increase in those living in poverty, the number of annual murders per million is expected to be lower by 29.91 on average.
- (d) For each percentage increase annual murders per million, the percentage of those living in poverty is expected to be higher by 2.56 on average.

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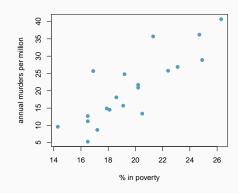
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Your turn

Suppose you want to predict annual murder count (per million) for a series of districts that were not included in the dataset. For which of the following districts would you be most comfortable with your prediction?

A district where % in poverty =

- (a) 5%
- (b) 15%
- (c) 20%
- (d) 26%
- (e) 40%

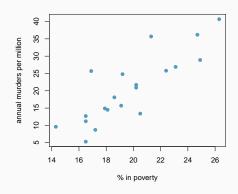


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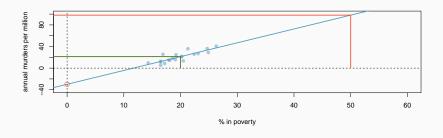
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A note about the intercept

Sometimes the intercept might be an extrapolation: useful for adjusting the height of the line, but meaningless in the context of the data.



Calculating predicted values

By hand:
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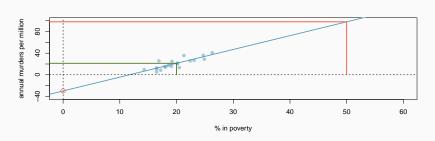
The predicted number of murders per million per year for a county with 20% poverty rate is:

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$$\widehat{murder} = -29.91 + 2.56 \times 20 = 21.29$$



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We can think of social science variables as comprised of two parts:

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Random variables contain both components

We can best understand random variables using probability distributions

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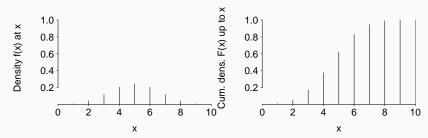
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Thus, for discrete distributions, the cdf is the <u>cumulative sum</u> of the pdf:

$$F(Y) = \sum_{\forall Y \le y} f(Y)$$

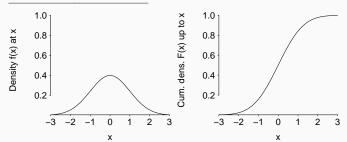
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Thus for continuous distributions, the cdf is the integral of the pdf:

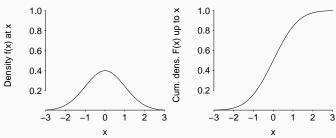
$$F(Y) = \int_{-\infty}^{y} f(Y) \mathrm{d}y$$

The Normal (Gaussian) distribution

$$f_{\mathcal{N}}(y|\mu,\sigma^2) = (2\pi\sigma^2)^{-1/2} \exp\left[\frac{-(y_i - \mu)^2}{2\sigma^2}\right]$$

Moments:
$$E(y) = \mu$$
 $Var = \sigma^2$

The Normal distribution is continuous and symmetric, with positive probability everywhere from $-\infty$ to ∞



The Normal distribution

What's the big deal about the Normal distribution?

One point of view: perhaps most continuous data are roughly Normally distributed

Why do people believe this?

They think the Central Limit Theorem applies to most data

The Central Limit Theorem

Suppose we have N independent random variables x_1, x_2, x_3, \ldots

Each x has an arbitrary probability distribution with mean μ_i and variance $\sigma_i^2 < \infty$

That is to say, these variables are not only independent, they could each have totally different distributions

Now suppose we average them all together into one super-variable,

$$X = \frac{1}{N} \sum_{i} x_{i}$$

The CLT shows that the distribution of this new variable, X, approaches a Normal distribution as $N \to \infty$

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With the Normal distribution in mind, recall the linear regression model

$$y_i = \beta_0 + \beta_1 x_i + \varepsilon_i$$

 $arepsilon_i$ is a normally distributed disturbance with mean 0 and variance σ^2

Equivalently, we write $\varepsilon_i \sim N(0, \sigma^2)$

Note that:

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The errors are assumed uncorrelated: $E(\varepsilon_i \times \varepsilon_j) = 0$ for all $i \neq j$

Recalling the definition of variance, note that in linear regression:

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$$\sigma^{2} = E\left((\varepsilon - E(\varepsilon))^{2}\right)$$
$$= E\left((\varepsilon - 0)^{2}\right)$$
$$= E(\varepsilon^{2})$$

The square root of σ^2 is known as the standard error of the regression

It is how much we expect y to differ from its expected value, $\beta_0 + \beta_1 x_i$, on average

Scalar representation:

$$y_i = \beta_0 + \beta_1 x_{1i} + \beta_2 x_{2i} + \varepsilon_i$$

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$$egin{array}{llll} \mathbf{y} & = & \mathbf{X} & oldsymbol{eta} & + & arepsilon \ n imes 1 & n imes k & k imes 1 & n imes 1 \end{array}$$

Writing out the matrices:

$$\begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{bmatrix} = \begin{bmatrix} 1 & x_{11} & x_{21} & \dots & x_{k1} \\ 1 & x_{12} & x_{22} & \dots & x_{k2} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 1 & x_{1n} & x_{2n} & \dots & x_{kn} \end{bmatrix} \begin{bmatrix} \beta_0 \\ \beta_1 \\ \vdots \\ \beta_k \end{bmatrix} + \begin{bmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \vdots \\ \varepsilon_n \end{bmatrix}$$

Note that we now have a vector of disturbances.

They have the same properties as before, but we will write them in matrix form.

The disturbances are still mean zero.

$$E(\varepsilon) = \begin{bmatrix} E(\varepsilon_1) \\ E(\varepsilon_2) \\ \vdots \\ E(\varepsilon_n) \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ \vdots \\ 0 \end{bmatrix}$$

But now we have an entire matrix of variances and covariances, Σ

$$\Sigma = \begin{bmatrix} \operatorname{var}(\varepsilon_1) & \operatorname{cov}(\varepsilon_1, \varepsilon_2) & \dots & \operatorname{cov}(\varepsilon_1, \varepsilon_n) \\ \operatorname{cov}(\varepsilon_2, \varepsilon_1) & \operatorname{var}(\varepsilon_2) & \dots & \operatorname{cov}(\varepsilon_2, \varepsilon_n) \\ \vdots & \vdots & \ddots & \vdots \\ \operatorname{cov}(\varepsilon_n, \varepsilon_1) & \operatorname{cov}(\varepsilon_n, \varepsilon_2) & \dots & \operatorname{var}(\varepsilon_n) \end{bmatrix}$$

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$$= \begin{bmatrix} \operatorname{E}(\varepsilon_{1}^{2}) & \operatorname{E}(\varepsilon_{1}\varepsilon_{2}) & \dots & \operatorname{E}(\varepsilon_{1}\varepsilon_{n}) \\ \operatorname{E}(\varepsilon_{2}\varepsilon_{1}) & \operatorname{E}(\varepsilon_{2}^{2}) & \dots & \operatorname{E}(\varepsilon_{2}\varepsilon_{n}) \\ \vdots & \vdots & \ddots & \vdots \\ \operatorname{E}(\varepsilon_{n}\varepsilon_{1}) & \operatorname{E}(\varepsilon_{n}\varepsilon_{2}) & \dots & \operatorname{E}(\varepsilon_{n}^{2}) \end{bmatrix}$$

However, the above matrix can be written far more compactly as an outer product

$$\boldsymbol{\Sigma} = \boldsymbol{\varepsilon} \boldsymbol{\varepsilon}'$$

Recall $\mathrm{E}(\varepsilon_i \varepsilon_j) = 0$ for all $i \neq j$, so all of the off-diagonal elements above are zero by assumption Recall also that all ε_i are assumed to have the same variance, σ^2 So if the linear regression assumptions hold, the variance-covariance matrix has a simple form:

$$\Sigma = \begin{bmatrix} \sigma^2 & 0 & \dots & 0 \\ 0 & \sigma^2 & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & \sigma^2 \end{bmatrix} = \sigma^2 \mathbf{I}$$

When these assumptions do not hold, we will need more complex models than simple linear regression

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The sum of squared errors can be written out as

$$arepsilon'arepsilon = (\mathsf{y} - \mathsf{X}eta)'(\mathsf{y} - \mathsf{X}eta)$$

(what is this notation doing? why do we need the transpose?)

We need two bits of matrix algebra:

$$(\mathbf{A} + \mathbf{B})' = \mathbf{A}' + \mathbf{B}'$$

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Simplifying, we get

$$\varepsilon' \varepsilon = \mathbf{y}' \mathbf{y} - 2\beta' \mathbf{X}' \mathbf{y} + \beta' \mathbf{X}' \mathbf{X} \beta$$

Now we need to take the derivative with respect to β , to see which β minimize the sum of squares.

How do we take the derivative of a scalar with respect to a vector? It's just a bunch of scalar derivatives stacked together:

$$\frac{\partial y}{\partial \mathbf{x}} = \begin{bmatrix} \frac{\partial y}{\partial x_1} & \frac{\partial y}{\partial x_2} & \dots & \frac{\partial y}{\partial x_n} \end{bmatrix}'$$

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For example, for **a** and **x** both $n \times 1$ vectors

$$y = \mathbf{a}'\mathbf{x} = a_1x_1 + a_2x_2 + \dots + a_nx_n$$
$$\frac{\partial y}{\partial \mathbf{x}} = \begin{bmatrix} a_1 & a_2 & \dots & a_n \end{bmatrix}'$$
$$\frac{\partial y}{\partial \mathbf{x}} = \mathbf{a}$$

A similar pattern holds for quadratic expresssions.

Note the vector analogue of x^2 is the inner product $\mathbf{x}'\mathbf{x}$

And the vector analogue of ax^2 is $\mathbf{x}'\mathbf{A}\mathbf{x}$, where \mathbf{A} is an $n \times n$ matrix of coefficients

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$$\frac{\partial ax^2}{\partial x} = 2ax$$
$$\frac{\partial x' Ax}{\partial x} = 2Ax$$

The details are a bit more complicated ($\mathbf{x}'\mathbf{A}\mathbf{x}$ is the sum of a lot of terms), but the intuition is the same.

$$\varepsilon' \varepsilon = \mathbf{y}' \mathbf{y} - 2\beta' \mathbf{X}' \mathbf{y} + \beta' \mathbf{X}' \mathbf{X} \beta$$

Take the derivative of the expression, setting it = 0, we get

$$\frac{\partial \varepsilon' \varepsilon}{\partial \beta} = -2\mathbf{X}'\mathbf{y} + 2\mathbf{X}'\mathbf{X}\beta = 0$$

This is a minimum, and the eta's that solve this equation thus minimize the sum of squares.

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So let's solve for β :

$$X'X\beta = X'y$$

$$\hat{oldsymbol{eta}} = (\mathbf{X}'\mathbf{X})^{-1}\mathbf{X}'\mathbf{y}$$

This is the least squares estimator for $oldsymbol{eta}$

As long as we have software to help us with matrix inversion, it is easy to calculate.

Outline

1. Housekeeping

2. Main ideas

- 1. Correlation coefficient describes the strength and direction of the linear association between two numerical variables
 - 2. Least squares line minimizes squared residuals
 - 3. Interpreting the least squares line
 - 4. Predict, but don't extrapolate
 - Regression in social sciences
 - 6. Distributions
 - 7. Linear Regression in Matrix Form

3. Assumptions

Is $(\mathbf{X}'\mathbf{X})^{-1}\mathbf{X}'\mathbf{y}$ a good estimate of β ?

Would another estimator be better?

What would an alternative be?

Maybe minimizing the sum of absolute errors?

Or something nonlinear?

First we'll have to decide what makes an estimator good.

Some common criteria:

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Bias

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Although it seems "obvious" on face that we always prefer an unbiased estimator if one is available we also want the estimate to be close to the truth most of the time

Unbiased methods are not perfect.

They usually still miss the truth by some amount, But the direction in which they miss is not systematic or known ahead of time.

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Biased estimates are not <u>necessarily</u> terrible.

A biased estimate of the time of day: a clock that is 2 minutes fast.

Efficiency:

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Stopped clock.	No	No
Random clock.	Yes	No
Clock that is "a lot fast"	No	No
Clock that is "a little fast"	No	Yes
A well-run atomic clock	Yes	Yes

To measure efficiency, we use **mean squared error**:

MSE =
$$E\left[\left(\beta - \hat{\beta}\right)^2\right]$$

= $Var(\hat{\beta}) + Bias(\hat{\beta}|\beta)^2$

 $\sqrt{\textit{MSE}}$ is how much you miss the truth by on average

In most cases, we want to use the estimator that minimizes MSE We will be especially happy when this is also an unbiased estimator But it won't always be

Consistency:

Consistency: An estimator that converges to the truth as the number of observations grows

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$$\mathrm{E}(\hat{\beta}-\beta) \to 0$$
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We will be mainly concerned with efficiency, secondarily with bias, and hardly at all with consistency

Two things that can go wrong:

- omitted variable bias
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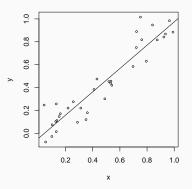
 Average children per marriage is 2.5. How many were in your family growing up? Are these numbers different? Who is "left out" in the second sample?

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- Average children per marriage is 2.5. How many were in your family growing up? Are these numbers different? Who is "left out" in the second sample?
- In testimony to NY state senate, motorcyclists testified that in their (multiple) crashes, helmets would not have prevented injuries. Who didn't testify?
- Regression example: Selection on the observed variables

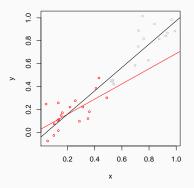
Selection bias



Suppose we conducted a survey & asked people their income (x) and conservatism (y)

With the full range of respondents, we find a strong relationship

Selection bias

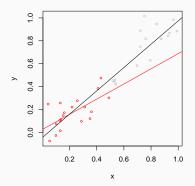


But suppose high income (or highly conservative) people decline to answer

Then we run a regression on the red dots only.

And get a result biased towards 0.

Selection bias



 \rightarrow Try to maximize variance of covariates, and avoid selecting on response variables

Most selection is unintentional, so think hard about sources of selection bias

What else can go wrong in a linear regression?

Even if your data are sampled without bias from the population of interest, and your model correctly specified, several data problems can violate the linear regression assumptions

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In order of declining severity:

Perfect collinearity

Endogeneity of covariates

Heteroskedasticity

Serial correlation

Non-normality

Lots of new jargon. Let's work through it.

Perfect Collinearity

Perfect collinearity occurs when $\mathbf{X}'\mathbf{X}$ is singular; ie, the determinant $|\mathbf{X}'\mathbf{X}|=0$

Happens when two or more columns of \boldsymbol{X} are linearly dependent on each other

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Very rare—except in panel data, as we will see

Matrix inversion—and thus LS regression—is impossible here

What if our covariates are correlated but not perfectly so?

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Those large se's are correct

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Have highly correlated **X** and large se's?

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Have highly correlated ${\bf X}$ and large se's? Then you lack sufficient data to precisely answer your research question

Exogenous & endogenous variables

So far, we have (implicitly) taken our regressors, X, as fixed

 \boldsymbol{X} is not dependent on \boldsymbol{Y}

 ${\sf Fixed} = {\sf pre-determined} = {\sf exogenous}$

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X is not dependent on Y

Fixed = pre-determined = exogenous

Y consists of a function of X plus an error

 ${f Y}$ is thus endogenous to ${f X}$

endogenous = "determined within the system"

Exogenous & endogenous variables

What if \boldsymbol{Y} helps determine \boldsymbol{X} in the first place?

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Very common in political science:

- campaign spending and share of the popular vote.
- policy attitudes and party identification
- arms races and war, etc.
- exchange rate policy and inflation

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In these cases, \boldsymbol{Y} and \boldsymbol{X} are both endogenous

Least squares is biased in this case

It will remain biased even as you add more data

In other words, it is inconsistent, or biased even as $N o \infty$

Linear regression allows us to model the mean of a variable well

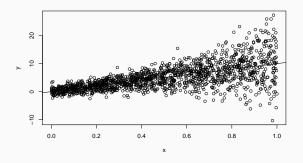
Y could be any linear function of β and X

But LS always assumes the variance of that variable is the same:

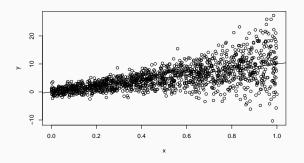
 σ^2 , a constant

We don't think **Y** has constant mean. Why expect constant variance?

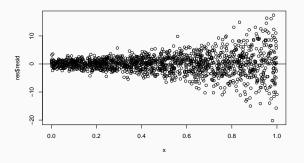
In fact, heteroskedasticity—non-constant error variance—is very common



A common pattern of heteroskedasticity: Variance and mean increase together Here, they are both correlated with the covariate X

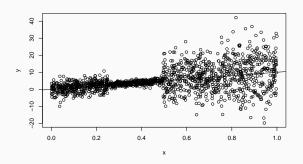


In a fuzzy sense, X is a necessary but not sufficient condition for Y. This is usually an important point substantively. Heteroskedasticity is interesting, not just a nuisance



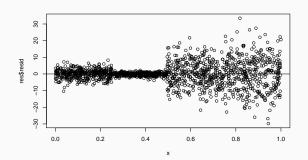
We can usually find heteroskedasticity by plotting the residuals against each covariate

Look for a pattern. Often a megaphone



But other patterns are possible.

Above, there is a dramatic difference in variance in different parts of the dataset



The same diagnostic reveals this problem.

Heteroskedasticity of this type often appears in panel datasets, where there are groups of observations from different units that each share a variance

Unpacking σ^2

Every observation consists of a systematic component $(\mathbf{x}_i\beta)$ and a stochastic component (ε_i)

Generally, we can think of the stochastic component as an n-vector ε following a multivariate normal distribution:

$$arepsilon \sim \mathcal{MVN}(\mathbf{0}, oldsymbol{\Sigma})$$

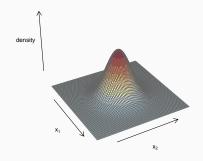
Aside: how the Multivariate Normal distribution works

Consider the simplest multivariate normal distribution, the joint distribution of two normal variables x_1 and x_2

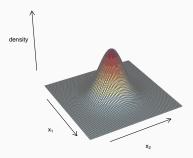
As usual, let μ indicate a mean, and σ a variance or covariance

$$egin{array}{lcl} egin{array}{lcl} eta_1 & eta_1 & \sigma_{1,2} \ egin{array}{lcl} egin{array}{lcl}$$

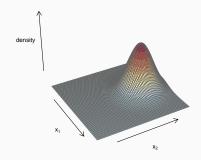
The MVN is more than the sum of its parts: There is a mean and variance for each variable, <u>and</u> covariance between each pair



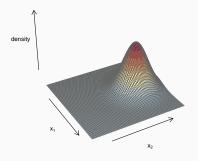
$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN}\left(\left[\begin{array}{c} \mathbf{0} \\ \mathbf{0} \end{array}\right], \left[\begin{array}{cc} \mathbf{1} & \mathbf{0} \\ \mathbf{0} & \mathbf{1} \end{array}\right]\right)$$



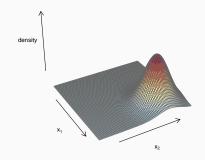
The standard MVN, with zero means, unit variances, and no covariance, looks like a higher dimension version of the normal: a symmetric mountain of probability



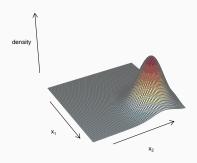
$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN}\left(\left[\begin{array}{c} 0 \\ 2 \end{array}\right], \left[\begin{array}{c} 1 & 0 \\ 0 & 1 \end{array}\right]\right)$$



Shifting the mean of \mathbf{x}_2 moves the MVN in one dimension only Mean shifts affect only one dimension at a time

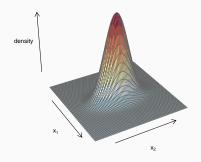


$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN}\left(\left[\begin{array}{c} 2 \\ 2 \end{array}\right], \left[\begin{array}{c} 1 & 0 \\ 0 & 1 \end{array}\right]\right)$$

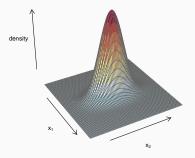


We could, of course, move the means of our variables at the same time.

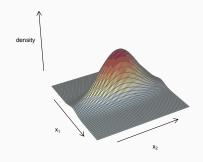
This MVN says the most likely outcome is both x_1 and x_2 will be near 2.0



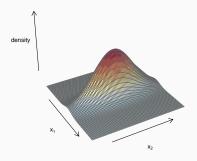
$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN}\left(\left[\begin{array}{c} 0 \\ 0 \end{array}\right], \left[\begin{array}{c} 0.33 & 0 \\ 0 & 1 \end{array}\right]\right)$$



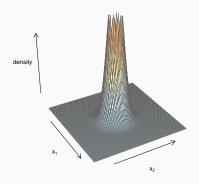
Shrinking the variance of x_1 moves the mass of probability towards the mean of x_1 , but leaves the distribution around x_2 untouched



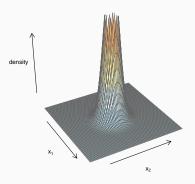
$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN} \left(\left[\begin{array}{c} 0 \\ 0 \end{array}\right], \left[\begin{array}{cc} 0.33 & 0 \\ 0 & 3 \end{array}\right] \right)$$



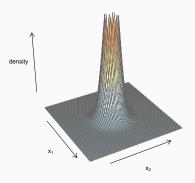
Increasing the variance of x_2 spreads the probability out, so we are less certain of x_2 , but just as certain of x_1 as before



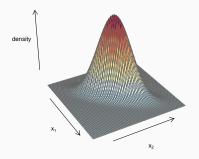
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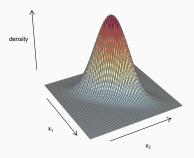
If the variance is small on all dimensions, the distribution collapses to a spike over the means of all variables



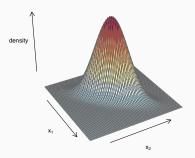
In this case, we are fairly certain of where all our variables tend to lie



$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN} \left(\left[\begin{array}{c} 0 \\ 0 \end{array}\right], \left[\begin{array}{cc} 1 & 0.8 \\ 0.8 & 1 \end{array}\right] \right)$$

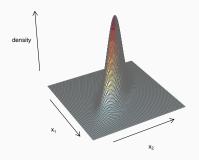


In this special case, with unit variances, the covariance is also the correlation, so our distribution say x_1 and x_2 are correlated at r=0.8

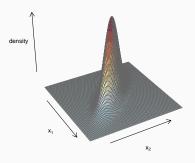


A positive correlation between our variables makes the $\ensuremath{\mathsf{MVN}}$ asymmetric,

with greater mass on likely combinations



$$\left[\begin{array}{c} \mathbf{x}_1 \\ \mathbf{x}_2 \end{array}\right] = \mathcal{MVN} \left(\left[\begin{array}{c} 0 \\ 0 \end{array}\right], \left[\begin{array}{cc} 1 & -0.8 \\ -0.8 & 1 \end{array}\right]\right)$$



A negative correlation makes $\underline{\text{mismatched}}$ values of our covariates more likely

In our current example, we have a huge multivariate normal distribtion:

each observation has its own mean and variance, and a covariance with every other observation

Suppose we have four observations. The Var-cov matrix of the disturbances is then

$$\Sigma = \begin{bmatrix} \sigma_1^2 & \sigma_{12} & \sigma_{13} & \sigma_{14} \\ \sigma_{21} & \sigma_2^2 & \sigma_{23} & \sigma_{24} \\ \sigma_{31} & \sigma_{32} & \sigma_3^2 & \sigma_{34} \\ \sigma_{41} & \sigma_{42} & \sigma_{43} & \sigma_4^2 \end{bmatrix}$$

In its most "ordinary" form, linear regression puts strict conditions on the variance-variance matrix, Σ

Again, assuming we have only four observations, the Var-cov matrix is

$$\mathbf{\Sigma} = \sigma^2 \mathbf{I} = \left[egin{array}{cccc} \sigma^2 & 0 & 0 & 0 \ 0 & \sigma^2 & 0 & 0 \ 0 & 0 & \sigma^2 & 0 \ 0 & 0 & 0 & \sigma^2 \end{array}
ight]$$

Could treat each observation as consisting of $\mathbf{x}_i \boldsymbol{\beta}$ and a separate, univariate normal disturbance, each with the same variance, σ^2 .

This is the usual linear regression set up

Will look like our first example MVN: a symmetric mountain, but

Suppose the distrurbances are heteroskedastic.

Now each observation has an error term drawn from a Normal with its own variance

$$\Sigma = \left[\begin{array}{cccc} \sigma_1^2 & 0 & 0 & 0 \\ 0 & \sigma_2^2 & 0 & 0 \\ 0 & 0 & \sigma_3^2 & 0 \\ 0 & 0 & 0 & \sigma_4^2 \end{array} \right]$$

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ight]$$

Still no covariance across disturbances.

Even so, we now have more parameters than we can estimate.

If every observation has its own unknown variance, we cannot estimate them

This MVN looks like the first example of a ridge: steeper in some directions than others, but not "tilted"

Heteroskedasticity does <u>not</u> bias least squares

But LS is inefficient in the presence of heteroskedasticity

More efficient estimators give greater weight to observations with low variance

They pay more attention to the signal, and less attention to the noise

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More efficient estimators give greater weight to observations with low variance

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Heteroskedasticity tends to make se's incorrect, because they depend on the estimate of σ^2

Researchers often try to "fix" standard errors to deal with this (more on this later)

Unpacking σ^2 : heteroskedasticity & autocorrelation

Suppose each disturbance has its own variance, and may be correlated with other disturbances

The most general case allows for both heteroskedasticity & autocorrelation

$$\Sigma = \begin{bmatrix} \sigma_1^2 & \sigma_{12} & \sigma_{13} & \sigma_{14} \\ \sigma_{21} & \sigma_2^2 & \sigma_{23} & \sigma_{24} \\ \sigma_{31} & \sigma_{32} & \sigma_3^2 & \sigma_{34} \\ \sigma_{41} & \sigma_{42} & \sigma_{43} & \sigma_4^2 \end{bmatrix}$$

LS is unbiased but inefficient in this case

The standard errors will be wrong, however

Key application: time series.

Current period is usually a function of the past

Gauss-Markov Conditions

So when is least squares unbiased?

When is it efficient?

When are the standard errors correct?

To judge the performance	of LS,	we'll need	to make	some ass	sumptions
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3	Disturbances have mean 0	$\mathrm{E}(arepsilon)=0$	Biased, even as $N o \infty$
4	No serial correlation	$\mathrm{E}(\varepsilon_i\varepsilon_j)=0, i\neq j$	Unbiased but ineff. se's wrong
5	Homoskedastic errors	$\mathrm{E}(oldsymbol{arepsilon}'oldsymbol{arepsilon}) = \sigma^2 \mathbf{I}$	

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5	Homoskedastic errors	$\mathrm{E}(arepsilon'arepsilon) = \sigma^2 \mathbf{I}$	Unbiased but ineff. se's wrong
6	Gaussian error distrib	$arepsilon \sim \mathcal{N}(0,\sigma^2)$	se's wrong unless $N o \infty$

(Assumptions get stronger from top to bottom, but 4 & 5 could be combined)

Gauss-Markov Theorem

It is easy to show β_{LS} is linear and unbiased, under Asps 1–3:

If
$$\mathbf{y} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\varepsilon}$$
, $\mathrm{E}(\boldsymbol{\varepsilon}) = 0$, then by substitution

$$\hat{eta}_{ ext{LS}} = (\mathbf{X}'\mathbf{X})^{-1}\mathbf{X}'(\mathbf{X}eta + arepsilon)$$

$$= eta + (\mathbf{X}'\mathbf{X})^{-1}\mathbf{X}'arepsilon$$

So long as

- (X'X)⁻¹ is uniquely identified,
- ullet X is exogenous or at least uncorrelated with arepsilon, and
- $E(\varepsilon) = 0$ (regardless of the distribution of ε)

Then
$$\mathbf{E}(\hat{oldsymbol{eta}}_{\mathrm{LS}}) = oldsymbol{eta}$$

 $ightarrow eta_{
m LS}$ is unbiased and a linear function of ${f y}$.

Gauss-Markov Theorem

If we make assumptions 1–5, we can make a stronger claim

When there is no serial correlation, no heteroskedasticity, no endogeneity, and no perfect collinearity, then

Gauss-Markov holds that LS is the best linear unbiased estimator (BLUE)

BLUE means that among linear estimators that are unbiased, $\hat{\beta}_{LS}$ has the least variance.

But, there might be a nonlinear estimator with lower MSE overall, unless . . .

If in addition to Asp 1-5, the disturbances are normally distributed (6), then

Gauss-Markov holds LS is Minimum Variance Unbiased (MVU)