Multiple Regression Analysis



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- We now return to the issue of changes in scale and origin we met before in Chapter 2 and examine the effects of rescaling the dependent or independent variables on *se*, *t* statistics, *F* statistics, and *CI*.
- As expected, when variables are rescaled, the coefficients, *se*, *CI*, *t* and *F* statistics change in ways that preserve all measured effects and testing outcomes.
- Hence, our conclusions are not affected by the units of measurement in the variables involved.

Consider the following estimated equation:

$$\hat{y} = \hat{\beta}_0 + \hat{\beta}_1 x_1 + \hat{\beta}_2 x_2$$

and now consider what happens to our OLS statistics as we change the scale and origin of y and of x_1 .



We can work out these effects by simply manipulating the above equation.

1. Changes in the scale of $y: c_1.y$

$$c_1.\hat{y} = (c_1.\hat{\beta}_0) + (c_1.\hat{\beta}_1)x_1 + (c_1.\hat{\beta}_2)x_2$$

✓ Coefficients are multiplied by c₁.
 ✓ Standard errors are multiplied by c₁.
 ✓ Statistical significance is not affected.
 ✓ *CI* change by the same factor, c₁.

Residuals are multiplied by c_1 . ✓ SSR are multiplied by c_1^2 . ✓ Standard Error of the Regression, SER = $\hat{\sigma}$, is multiplied by c_1 . \checkmark R^2 is not affected, so the overall significance of the regression is not affected.

2. Changes in the origin of *y*: $c_0 + y$

$$c_0 + \hat{y} = (c_0 + \hat{\beta}_0) + \hat{\beta}_1 x_1 + \hat{\beta}_2 x_2$$

- Only the intercept, β₀, is affected.
 The slope coefficients, measuring partial effects, are not affected.
- Residuals are not affected.
- \checkmark R^2 is not affected.

3. Changes in the scale of $x_1: d_1.x_1$

$$\hat{y} = \hat{\beta}_0 + (\hat{\beta}_1 / d_1)(d_1 \cdot x_1) + \hat{\beta}_2 x_2$$

The coefficient associated to x₁, β₁, is divided by d₁.
 All other coefficients are not affected.
 The standard error of β₁ is divided by d₁.
 Statistical significance is not affected.
 The *CI* for β₁ change by the factor, 1/d₁.

- Residuals are not affected.
- Hence, neither SSR nor the SER are affected.
- $\checkmark R^2 \text{ is not affected, so the overall} \\ \text{significance of the regression is not} \\ \text{affected.}$

4. Changes in the origin of $x_1: d_0 + x_1$

$$\hat{y} = (\hat{\beta}_0 - \hat{\beta}_1 d_0) + \hat{\beta}_1 (x_1 + d_0) + \hat{\beta}_2 x_2$$

- Only the intercept, β₀, is affected.
 The slope coefficients, measuring partial effects, are not affected.
- Residuals are not affected.
- \checkmark R^2 is not affected.

- Conclusion: Changes in scale and/or origin does not affect to any substantial part of the regression.
- In particular, statistical significance and interpretation of coefficients is not affected by data scaling.



Note that to make our equation invariant to the origin of the variables we need an intercept in our equation.

 This analysis shows clearly that if variables appear in logarithmic form, changing the units of measurement does not affect the slope coefficients.

This follows from the fact that $log(c_1, y) = log(c_1) + log(y)$ $c_1 > 0$

 $\log(d_1.x_j) = \log(d_1) + \log(x_j)$ $d_1 > 0$

so only the intercept is affected in these cases.

- Sometimes in econometric applications, a key variable is measured on a scale that is difficult to interpret, for example, test scores, synthetic indexes,...
- In such cases, we can be interested in see what happens to y when the corresponding independent variable varies by one standard deviation.

Sometimes, it is useful to obtain regression results when all variables involved, y as well as the x's, have been standardized.



To standardize a variable subtracts its mean and divide by its standard deviation.



Why is standardization useful?

Lets see what this transformation implies for the coefficient estimates.

Beta Coefficients $y_{i} = \hat{\beta}_{0} + \hat{\beta}_{1}x_{i1} + \hat{\beta}_{2}x_{i2} + \ldots + \hat{\beta}_{k}x_{ik} + \hat{u}_{i}$ Averaging this equation and subtracting $y_{i} - \overline{y} = \hat{\beta}_{1}(x_{i1} - \overline{x}_{1}) + \hat{\beta}_{2}(x_{i2} - \overline{x}_{2}) + \ldots + \hat{\beta}_{k}(x_{ik} - \overline{x}_{k}) + \hat{u}_{i}$

Simple algebra gives us the estimated equation in standardized form

$$\frac{y_i - \overline{y}}{\hat{\sigma}_y} = \left(\hat{\beta}_1 \frac{\hat{\sigma}_1}{\hat{\sigma}_y}\right) \left[\frac{(x_{i1} - \overline{x}_1)}{\hat{\sigma}_1}\right] + \left(\hat{\beta}_2 \frac{\hat{\sigma}_2}{\hat{\sigma}_y}\right) \left[\frac{(x_{i2} - \overline{x}_2)}{\hat{\sigma}_2}\right] + \dots + \left(\hat{\beta}_k \frac{\hat{\sigma}_k}{\hat{\sigma}_y}\right) \left[\frac{(x_{ik} - \overline{x}_k)}{\hat{\sigma}_k}\right] + \frac{\hat{u}_i}{\hat{\sigma}_y}$$

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Which we can rewrite as

$$z_y = \hat{b}_1 z_1 + \hat{b}_2 z_2 + \ldots + \hat{b}_k z_k + \hat{e}_1$$

where z denotes an standardized variable, the zscore, \hat{e} denotes the error and the new coefficients are

$$\hat{b}_j = \hat{\beta}_j \frac{\sigma_j}{\hat{\sigma}_v}$$
 for $j = 1, 2, ..., k$

These \hat{b}_j are traditionally called **standardized coefficients** or **beta coefficients**.

 \diamond The meaning of these coefficients is as follows: If x_i increases by one standard deviation, then \hat{y} changes by b_i standard deviations, holding all other variables constant.



Thus, we are measuring effects not in terms of the original units of y and x_i , but in standard deviation units.

Because the equation in terms of the z-score makes the scale of the regressors irrelevant, this equation puts the explanatory variables on equal footing.

♦ In a standard OLS equation, it is not possible to simply look at the size of different coefficients and conclude that the explanatory variable with the largest coefficient is "the most important".



• We just have seen that the magnitudes of coefficients can be changed at will by changing the scale of x_i .



 \diamond But, when each x_i has been standardized, comparing magnitudes of the resulting beta coefficients is more compelling.

Functional Form

- OLS can be used for modeling relationships that are not strictly linear in x and y by using nonlinear functions of x and y, if the model is still linear in the parameters.
- We consider some possibilities that often appear in applied work:
- 1. $\log' s \text{ of } x \text{ and } y$.
- 2. quadratic forms of *x*.
- 3. Interactions of *x* variables.

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Proportions and Percentages



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Proportions and Percentages

4. Changes in logarithms:



Hence,

$100.\Delta \log(x) \approx \%.\Delta x$

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A Linear Model for log(y)Consider the model $\log(y) = \beta_0 + \beta_1 x + u$ \diamond What is the meaning of β_1 in this model? • If $\Delta u = 0$, then x has a linear effect on log(y): $\Delta \log(y) = \beta_1 \Delta x$ or, $\%\Delta y = (100.\beta_1).\Delta x$ i.e. $100.\beta_1$ is the percentage change in y by unit of x.

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A Constant Elasticity Model

Consider the model $\log(y) = \beta_0 + \beta_1 \log(x) + u$ • What is the meaning of β_1 in this model? $If \Delta u = 0$, then $\log(x)$ has a linear effect on log(y): $\Delta \log(y) = \beta_1 \Delta \log(x) \iff \% \Delta y = \beta_1.\% \Delta x$ i.e. β_1 is the elasticity of y with respect to x.

Functional Forms Involving logs

Model	Dependent	Independent	Interpretation
	Variable	Variable	of β_1
level-level	у	X	$\Delta y = \beta_1 . \Delta x$
level-log	У	$\log(x)$	$\Delta y = (\beta_1 / 100).\% \Delta x$
log-level	$\log(y)$	X	$\% \Delta y = (100.\beta_1).\Delta x$
log-log	$\log(y)$	$\log(x)$	$\% \Delta y = \beta_1.\% \Delta x$

Functional Form

Important: While the mechanics of the linear regression does not depend on how y and the x's are defined, the interpretation of the coefficients does depend on their definitions.

Why use log models?

- Using log's leads to coefficients with appealing interpretations, i.e. elasticity or semi-elasticity. ✓ Models with log's are invariant to the scale of the variables, since they measure proportional changes. \checkmark For models with y > 0, using $\log(y)$ as the dependent variable often satisfy the CLM assumptions more closely than models using the level of y. \checkmark For models with y > 0, the conditional distribution is
 - often heteroskedastic or skewed, while log(y) is much less so.

Why use log models?

- Taking log's usually narrows the range of the variable. This makes estimates less sensitive to outlying (or extreme) observations on the dependent or independent variables.
- One limitation of the log is that it can not be used if a variables can take zero or negative values.
- One drawback to using a dependent variable in log form is that it is more difficult to predict the original variable. The original model allows us to predict log(y), not y.

Why use log models?

- Also it is *not* legitimate to compare R² from models where y is the dependent variable in one case and log(y) is the dependent variable in the other. These measures explained variations in different variables.
- Important: This is a general rule, the R² cannot be used to compare models with different dependent variable.

Some Rules of Thumb

- What types of variables are often used in log form?
- \checkmark Variables in money terms that must be positive.
- Very large variables, such as population.
- What types of variables are often used in level form?
- ✓ Variables measured in years.
- Variables that are a proportion or percent, i.e. inflation, interest rates.

Quadratic Models

A quadratic model is of the form $y = \beta_0 + \beta_1 x + \beta_2 x^2 + u$

Quadratic functions are also used quite often in applied economics to capture decreasing or increasing marginal effects.
 Important: β₁ does not measure the change in *y* with respect to *x*; it makes no sense to hold x² fixed while changing *x*.

Quadratic Models

• If $\Delta u = 0$ then,

 $\Delta y \approx (\beta_1 + 2\beta_2 x) \Delta x \implies \frac{\Delta y}{\Lambda x} \approx \beta_1 + 2\beta_2 x$

the marginal effect of *x* on *y* depends linearly on the value of *x*.

The estimated slope is $\beta_1 + 2\beta_2 x$.

• In a particular application this marginal effect should be evaluated at interesting values of x.

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More on Quadratic Models

• Suppose that $\beta_1 > 0$ and $\beta_2 < 0$.

 \diamond Then y is increasing in x at first, but will eventually turn around and be decreasing in

 \boldsymbol{X} .

The turning point will be at

$$x^* = \left| \frac{\beta_1}{2\beta_2} \right|$$

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More on Quadratic Models

• Suppose that $\beta_1 < 0$ and $\beta_2 > 0$.

Then y is decreasing in x at first, but willeventually turn around and be increasing in

Х.

The turning point will be at

$x^* = \left| \frac{\beta_1}{2\beta_2} \right|$

which is the same as before.

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Interaction Terms

Sometimes, it is natural for the partial effect, elasticity or semi-elasticity of the dependent variable with respect to an explanatory variable to depend on the magnitude of yet another explanatory variable.

These effects can be modeled through interaction terms, $x_i x_j$.



Interaction Terms

- The partial effect of x_1 on y depends linearly on x_2 .
- In summarizing the effect of x₁ on y, we must evaluate the above expression at interesting and representative values of x₂, for examples the sample mean of x₂.

Functional Form

This shows clearly that the partial effects of x_j on y are constant only if the model is linear in variables. In all other cases the interpretation of the coefficients does depend on the definitions of the variables.

R-Squared

- We found before the *R*² as a **goodness of fit** measure.
- R^2 is simply an estimate of how much variation in y is explained by the x's, and even it is intuitively obvious that a higher R^2 is preferable to a lower one, nothing about the classical model assumptions requires that R^2 be above any particular value.
- A small R^2 does imply that the error variance is large relative to the variance of *y*, which means that the β_i are not precisely estimated.

R-Squared

- But remember, that a large error variance can be offset by a large sample size, so if *n* is large enough, we may be able to precisely estimate the partial effects even though we have not controlled for many unobserved factors.
- Also that the relative *change* in the R^2 , when variables are added to an equation, is very useful: the *F* statistic for testing the joint significance of the added variables crucially depends on the difference in the R^2 between the unrestricted and the restricted models.

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- Recall that the R^2 will always increase as more variables are added to a given model.
- This can lead to the false impression that models with more explanatory variables are always preferred, but this is completely false. If we add variables to a given model, R^2 will never decrease, even if these variables are not significant.



 \clubsuit To avoid this algebraic fact we can "adjust" the R^2 in a way that takes into account the number of variables included in a given the model.

 \clubsuit To see how the usual R^2 might be adjusted, it is usefully written as





 \clubsuit This expression reveals what R^2 is actually estimating.

• The population R^2 is defined as $1 - \frac{\sigma_u^2}{\sigma^2}$

This is what R² is supposed to be estimating.
 However, we have better estimates for these variances that the ones used in the R². So lets use unbiased estimates for these variances

$$\overline{R}^{2} = 1 - \frac{\text{SSR}/(n-k-1)}{\text{SST}/(n-1)} = 1 - (1 - R^{2}) \frac{n-1}{n-k-1}$$

\diamond This is the *adjusted* R^2 .

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 \checkmark The primary attractiveness of \overline{R}^2 is that it imposes a penalty for adding additional independent variables to a model. ♦ If an independent variable is added to a model then SSR falls, but so does the df in the regression, n - k - 1. So \overline{R}^2 can go up or down when a new independent variable is added to a regression.

An interesting algebraic fact is that if we add a new independent variable to a regression equation, \overline{R}^2 increases if, an only if, the *t* statistic on the new variable is greater than one in absolute value.

Thus we see immediately that using \overline{R}^2 to decide whether a certain independent variable belongs in a model gives us a different answer than standard *t* testing.

 \clubsuit It is important not to focus too much on \mathbb{R}^2 or \mathbb{R}^2 , and lose insights from economic theory and common sense.



• Goodness of fit by itself is not an objective.

♦ If economic theory clearly predicts a variable belongs to a model, generally leave it in.



On't try to include a variable that prohibits a sensible interpretation of the variables of interest. Remember the ceteris paribus interpretation of multiple regression.

- Provided the above conditions are fulfilled,
 you can use the R² to measure the goodness
 of fit of models with the same number of
 independent variables and the same y:
 - (1) $y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + u$ (2) $y = \beta_0 + \beta_1 x_1 + \beta_3 x_3 + u$

These are nonnested models, because neither equation is a special case of the other.

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 \clubsuit You can use the \overline{R}^2 to measure the goodness of fit of models with different number of independent variables **and** the same y:

(1) $y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \beta_3 x_3 + \beta_4 x_4 + u$ $y = \beta_0 + \beta_1 x_1 + \beta_3 x_3 + \beta_4 \log(x_4) + u$ (2)



Explanatory variables can appear with different functional form, but not y.

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- You **cannot** use neither the R^2 nor \overline{R}^2 to measure the goodness of fit of models with different functional forms for the dependent variable, y:
 - (1) $y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \beta_3 x_3 + \beta_4 x_4 + u$ (2) $\log(y) = \beta_0 + \beta_1 x_1 + \beta_4 \log(x_4) + u$

The reason is simple: the variation to be explained, SST, is different for both models.

Suppose we have estimated the equation

 $\hat{y} = \hat{\beta}_0 + \hat{\beta}_1 x_1 + \hat{\beta}_2 x_2 + \ldots + \hat{\beta}_k x_k$

 \diamond When we plug in particular values of the x's, we obtain a prediction for y, which is an estimate of the expected value of y given the particular values for the x's.



 \diamond Let c_1, c_2, \ldots, c_k denote the particular values for each of the k independent variables; these may or may no correspond to an actual data point in our sample.





large n-k-1 then $t_{.025} \approx 1.96$

- How do we obtain the *se* of $\hat{\theta}_0$?
- If the computer software does not do the job for you, note that all you need is a *se* of a linear combination of the OLS estimators, just as in hypothesis testing, so the same trick we used there works here.
- Write $\beta_0 = \theta_0 \beta_1 c_1 \beta_2 c_2 \dots \beta_k c_k$, and plug this into the equation

$$y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \ldots + \beta_k x_k + u$$

to obtain

 $y = \theta_0 + \beta_1 (x_1 - c_1) + \beta_2 (x_2 - c_2) + \ldots + \beta_k (x_k - c_k) + u$

- ♦ In other words, we subtract the value c_j from each observation on x_j, and then we run the regression of y_i on (x_{i1} c₁), (x_{i2} c₂), ..., (x_{ik} c_k), i = 1,...,n
 ♦ The predicted value, and more importantly, its se, are obtained from the *intercept*, or constant, in this regression.
- This result is not surprising, since intuitively we have less uncertainty near the middle of our data.

Prediction: CI



This illustrates graphically the confidence interval for predictions in the SLR case.

- The previous method allows us to put a *CI* around the OLS estimate of $E(y|x_1, x_2, ..., x_3)$, for any values of the x's.
- In other words, we obtain a *CI* for the average value of *y* for the subpopulation with a given set of covariates.
 - But a *CI* for the average unit in the subpopulation is not exactly the same as a *CI* for a particular unit in the subpopulation.
- In forming a *CI* for an unknown outcome on *y*, we must account for another very important source of variation: the variance in the unobserved error, which measures our ignorance on the unobserved factors that affect *y*.

 \clubsuit Let y^0 denote the value for which we would like to construct a CI, usually called prediction interval. Let $x_1^0, x_2^0, ..., x_k^0$ be the new values of the x's, which we observe, and let u^0 be the unobserved error. Therefore, we have

 $y^{0} = \beta_{0} + \beta_{1}x_{1}^{0} + \beta_{2}x_{2}^{0} + \dots + \beta_{k}x_{k}^{0} + u^{0}$



 \diamond As before, our best point prediction of y^0 is the expected value of y^0 given the explanatory variables, which we estimate from the OLS regression line $\hat{y}^{0} = \hat{\beta}_{0} + \hat{\beta}_{1}x_{1}^{0} + \hat{\beta}_{2}x_{2}^{0} + \dots + \hat{\beta}_{k}x_{k}^{0}$

- The **prediction error** in using \hat{y}^0 to predict y^0 is $\hat{e}^{0} = y^{0} - \hat{y}^{0} = (\beta_{0} + \beta_{1}x_{1}^{0} + \beta_{2}x_{2}^{0} + \dots + \beta_{k}x_{k}^{0}) + u^{0} - \hat{y}^{0}$ Because OLS estimators are unbiased and $E(u^0) = 0$, then $E(\hat{e}^0) = 0$. So the expected prediction error is zero.
- In finding the variance of \hat{e}^0 , note that u^0 is uncorrelated with \hat{y}^0 (why?).
- Therefore, the variance of the prediction error (conditional on the x's) is the sum of the variances $Var(\hat{e}^{0}) = Var(\hat{y}^{0}) + Var(u^{0}) = Var(\hat{y}^{0}) + \sigma^{2}$

- There are two sources of variation in \hat{e}^0 .
- 1. The sampling error in \hat{y}^0 , which arises because we have estimated the β_i .
- 2. The ignorance of the unobserved factors that affect y, which is reflected in σ^2 .

• Under the CLM assumptions \hat{e}^0 is also normally distributed (conditional on the x's). And using unbiased estimators of $Var(\hat{y}^0)$ and σ^2 , we can define the *se* of \hat{e}^0 as

$$se(\hat{e}^0) = \left\{ \left[se(\hat{y}^0) \right]^2 + \hat{\sigma}^2 \right\}^{\overline{2}}$$

• Using the same reasoning for the *t* statistic of the β_j ,



 $\frac{\hat{e}^0}{se(\hat{e}^0)}$ has a *t* distribution with *n*-*k*-1 *df*. Therefore,

 $\Pr\left[-t_{.025} \le \frac{\hat{e}^0}{se(\hat{e}^0)} \le t_{.025}\right] = .95$

where $t_{.025}$ is the 97.5th percentile in the t_{n-k-1} distribution.

• Plugging in $\hat{e}^0 = y^0 - \hat{y}^0$ and rearranging gives a 95% **prediction interval** for y^0 : $\hat{y}^0 \pm t_{.025}.se(\hat{e}^0)$.



 \diamond Usually the estimate of σ^2 is much larger than the variance of the prediction.



Thus, this prediction interval will be much wider than the simple *CI* for the prediction.

 \diamond As before with a large *df*, we can construct a 95% prediction interval using the *rule of thumb* $\hat{y}^0 \pm 2.se(\hat{e}^0)$, since for large n-k-1 then $t_{.025} \approx 1.96$.

Residual Analysis

- Sometimes, it is useful to examine the residuals for the individual observations. This process is known as residual analysis.
- Big residuals, either positive or negative, can be informative about special events or characteristics of individual observations.



Extreme residuals, greater in absolute value than 3 standard error of the regression, are called outliers. Outliers merit some consideration since they can

influence estimation results.

 \bigcirc Define *logy* = log(*y*), and consider the problem of predicting y when the estimated model is $logy = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \ldots + \beta_k x_k + u$ • Given OLS estimators we predict *logy* as $l\hat{o}gy = \hat{\beta}_0 + \hat{\beta}_1 x_1 + \hat{\beta}_2 x_2 + \ldots + \hat{\beta}_k x_k$ \diamond Simple exponentiation, $\hat{y} = \exp(l\hat{o}gy)$, will systematically *underestimate* the expected value of У. ♦ Instead, we need to scale this up by an estimate of the expected value of exp(u).

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Note that if $u \sim N(0, \sigma^2)$, then $E(\exp(u)) = \exp(\frac{\sigma^2}{2})$

Under the CLM assumptions MLR.1 through MLR.6, then $E(y \mid \boldsymbol{x}) = \exp(\sigma^2 / 2) \cdot \exp(\beta_0 + \beta_1 x_1 + \beta_2 x_2 + \ldots + \beta_k x_k)$ This equation shows that, under normality, the simple adjustment needed to predict y is $\hat{y} = \exp(\hat{\sigma}^2/2).\exp(l\hat{\sigma}gy)$ where $\hat{\sigma}^2$ is the unbiased estimator of σ^2 . \diamond Because $\hat{\sigma}^2 > 0 \implies \exp(\hat{\sigma}^2/2) > 1$



Given an estimate $\hat{\alpha}_0$, we can predict *y* as

 $\hat{y} = \hat{\alpha}_0 . \exp(l \hat{o} g y)$

- \diamond
- It turns out that a consistent estimator of $\hat{\alpha}_0$ is easily obtained:
- . Obtain the fitted values $l \hat{o} g y_i$
- 2. Create $\hat{m}_i = \exp(l\hat{o}gy_i)$
- 3. Regress y on \hat{m} , without an intercept. The coefficient on \hat{m} , the only coefficient there is, is the estimate of α_0 , i.e. $E(\exp(u))$.

4. Once $\hat{\alpha}_0$ is obtained, predict y as $\hat{y} = \hat{\alpha}_0 . \exp(l\hat{o}gy)$. Francisco J. Goerlich Introductory Econometrics 64

Comparing log(y) and y models

As mentioned before, R^2 cannot be used to compare models with different dependent variables. In particular, it cannot be used to compare models with y and log(y) as dependent variables.



♦ If the goal is to find a goodness-of-fit measure in the log(y) model that can be compared with the R^2 from a model where y is the dependent variable we can use the previous results.

 \clubsuit After running the regression of y on \hat{m} through the origin, we obtain the fitted values for this regression, $\hat{y}_i = \hat{\alpha}_0 \cdot \hat{m}_i$

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Comparing log(y) and y models



 \diamond Then, we find the sample correlation between \hat{y}_i and the actual y_i in the sample.

 \clubsuit The square of this can be compared with the R^2 we get by using y as the dependent variable in a linear regression model.

 \diamond Remember that the R^2 in the fitted equation $\hat{\mathbf{y}} = \hat{\beta}_0 + \hat{\beta}_1 x_1 + \hat{\beta}_2 x_2 + \ldots + \hat{\beta}_k x_k$ is just the squared correlation between y_i and \hat{y}_i .