## MULTIPLE REGRESSION BASICS

Documents prepared for use in course B01.1305, New York University, Stern School of Business

#### Introductory thoughts about multiple regression

page 3

Why do we do a multiple regression? What do we expect to learn from it? What is the multiple regression model? How can we sort out all the notation?

#### Scaling and transforming variables

page 9

Some variables cannot be used in their original forms. The most common strategy is taking logarithms, but sometimes ratios are used. The "gross size" concept is noted.

#### Data cleaning

page 11

Here are some strategies for checking a data set for coding errors.

#### Interpretation of coefficients in multiple regression

page 13

The interpretations are more complicated than in a simple regression. Also, we need to think about interpretations after logarithms have been used.

#### Pathologies in interpreting regression coefficients

page 15

Just when you thought you knew what regression coefficients meant . . .

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#### Regression analysis of variance table

page 18

Here is the layout of the analysis of variance table associated with regression. There is some simple structure to this table. Several of the important quantities associated with the regression are obtained directly from the analysis of variance table.

#### Indicator variables

page 20

Special techniques are needed in dealing with non-ordinal categorical independent variables with three or more values. A few comments relate to model selection, the topic of another document.

#### Noise in a regression

page 32

Random noise obscures the exact relationship between the dependent and independent variables. Here are pictures showing the consequences of increasing noise standard deviation. There is a technical discussion of the consequences of measurement noise in an independent variable. This entire discussion is done for simple regression, but the ideas carry over in a complicated way to multiple regression.

Cover photo: Praying mantis, 2003

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#### INPUT TO A REGRESSION PROBLEM

Simple regression:  $(x_1, Y_1), (x_1, Y_2), \dots, (x_n, Y_n)$ 

Multiple regression:  $((x1)_1, (x2)_1, (x3)_1, \dots (xK)_1, Y_1),$ 

 $((x1)_2, (x2)_2, (x3)_2, \dots (xK)_2, Y_2),$  $((x1)_3, (x2)_3, (x3)_3, \dots (xK)_3, Y_3),$ 

...,  $((x1)_n, (x2)_n, (x3)_n, ..., (xK)_n, Y_n),$ 

The variable *Y* is designated as the "dependent variable." The only distinction between the two situations above is whether there is just one *x* predictor or many. The predictors are called "independent variables."

There is a certain awkwardness about giving generic names for the independent variables in the multiple regression case. In this notation, x1 is the name of the first independent variable, and its values are  $(x1)_1, (x1)_2, (x1)_3, \ldots, (x1)_n$ . In any application, this awkwardness disappears, as the independent variables will have application-based names such as SALES, STAFF, RESERVE, BACKLOG, and so on. Then SALES would be the first independent variable, and its values would be  $SALES_1, SALES_2, SALES_3, \ldots, SALES_n$ .

The listing for the multiple regression case suggests that the data are found in a spreadsheet. In application programs like Minitab, the variables can appear in any of the spreadsheet columns. The dependent variable and the independent variables may appear in any columns in any order. Microsoft's EXCEL requires that you identify the independent variables by blocking off a section of the spreadsheet; this means that the independent variables must appear in consecutive columns.

#### MINDLESS COMPUTATIONAL POINT OF VIEW

The output from a regression exercise is a "fitted regression model."

Simple regression:  $\hat{Y} = b_0 + b_1 x$ 

Multiple regression:  $\hat{Y} = b_0 + b_1(x1) + b_2(x2) + b_3(x3) + ... + b_K(xK)$ 

Many statistical summaries are also produced. These are  $R^2$ , standard error of estimate, t statistics for the b's, an F statistic for the whole regression, leverage values, path coefficients, and on and on and on and ...... This work is generally done by a computer program, and we'll give a separate document listing and explaining the output.

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#### WHY DO PEOPLE DO REGRESSIONS?

A cheap answer is that they want to explore the relationships among the variables.

A slightly better answer is that we would like to use the framework of the methodology to get a yes-or-no answer to this question: Is there a significant relationship between variable *Y* and one or more of the predictors? Be aware that the word *significant* has a very special jargon meaning.

An simple but honest answer pleads curiousity.

The most valuable (and correct) use of regression is in making predictions; see the next point. Only a small minority of regression exercises end up by making a prediction, however.

#### HOW DO WE USE REGRESSIONS TO MAKE PREDICTIONS?

The prediction situation is one in which we have new predictor variables but do not yet have the corresponding *Y*.

Simple regression: We have a new x value, call it  $x_{new}$ , and the predicted (or

fitted) value for the corresponding Y value is

 $\hat{Y}_{\text{new}} = b_0 + b_1 x_{new}.$ 

Multiple regression: We have new predictors, call them  $(x1)_{new}$ ,  $(x2)_{new}$ ,  $(x3)_{new}$ ,

...,  $(xK)_{new}$ . The predicted (or fitted) value for the

corresponding Y value is

 $\hat{Y}_{new} = b_0 + b_1(x1)_{new} + b_2(x2)_{new} + b_3(x3)_{new} + \dots + b_K(xK)_{new}$ 

## CAN I PERFORM REGRESSIONS WITHOUT ANY UNDERSTANDING OF THE UNDERLYING MODEL AND WHAT THE OUTPUT MEANS?

Yes, many people do. In fact, we'll be able to come up with rote directions that will work in the great majority of cases. Of course, these rote directions will sometimes mislead you. And wisdom still works better than ignorance.

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#### WHAT'S THE REGRESSION MODEL?

The model says that Y is a linear function of the predictors, plus statistical noise.

Simple regression:  $Y_i = \beta_0 + \beta_1 x_i + \varepsilon_i$ 

Multiple regression:  $Y_i = \beta_0 + \beta_1 (x1)_i + \beta_2 (x2)_i + \beta_3 (x3)_i + ... + \beta_K (xK)_i + \varepsilon_i$ 

The coefficients (the  $\beta$ 's) are nonrandom but unknown quantities. The noise terms  $\epsilon_1$ ,  $\epsilon_2$ ,  $\epsilon_3$ , ...,  $\epsilon_n$  are random and unobserved. Moreover, we assume that these  $\epsilon$ 's are statistically independent, each with mean 0 and (unknown) standard deviation  $\sigma$ .

The model is simple, except for the details about the  $\varepsilon$ 's. We're just saying that each data point is obscured by noise of unknown magnitude. We assume that the noise terms are not out to deceive us by lining up in perverse ways, and this is accomplished by making the noise terms independent.

Sometimes we also assume that the noise terms are taken from normal populations, but this assumption is rarely crucial.

WHO GIVES ANYONE THE RIGHT TO MAKE A REGRESSION MODEL? DOES THIS MEAN THAT WE CAN JUST SAY SOMETHING AND IT AUTOMATICALLY IS CONSIDERED AS TRUE?

Good questions. Merely claiming that a model is correct does not make it correct. A model is a mathematical abstraction of reality. Models are selected on the basis of simplicity and credibility. The regression model used here has proved very effective. A careful user of regression will make a number of checks to determine if the regression model is believable. If the model is not believable, remedial action must be taken.

HOW CAN WE TELL IF A REGRESSION MODEL IS BELIEVABLE? AND WHAT'S THIS REMEDIAL ACTION STUFF?

Patience, please. It helps to examine some successful regression exercises before moving on to these questions.

# THERE SEEMS TO BE SOME PARALLEL STRUCTURE INVOLVING THE MODEL AND THE FITTED MODEL.

It helps to see these things side-by-side.

Simple regression:

The model is  $Y_i = \beta_0 + \beta_1 x_i + \varepsilon_i$ 

The <u>fitted model</u> is  $\hat{Y} = b_0 + b_1 x$ 

Multiple regression:

The model is  $Y_i = \beta_0 + \beta_1 (x1)_i + \beta_2 (x2)_i + \beta_3 (x3)_i + ... + \beta_K (xK)_i + \varepsilon_i$ 

The fitted model is  $\hat{Y} = b_0 + b_1(x1) + b_2(x2) + b_3(x3) + ... + b_K(xK)$ 

The Roman letters (the b's) are estimates of the corresponding Greek letters (the  $\beta$ 's).

#### WHAT ARE THE FITTED VALUES?

In any regression, we can "predict" or retro-fit the *Y* values that we've already observed, in the spirit of the PREDICTIONS section above.

Simple regression:

$$Y_i = \alpha + \beta x_i + \varepsilon_i$$

$$\hat{Y} = a + bx$$

The fitted value for point *i* is

$$\hat{Y}_i = a + bx_i$$

Multiple regression:

$$Y_i = \beta_0 + \beta_1 (x1)_i + \beta_2 (x2)_i + \beta_3 (x3)_i + \dots + \beta_K (xK)_i + \varepsilon_i$$

$$\hat{Y} = b_0 + b_1(x1) + b_2(x2) + b_3(x3) + ... + b_K(xK)$$

The fitted value for point i is

$$\hat{Y}_i = b_0 + b_1(x1)_i + b_2(x2)_i + b_3(x3)_i + ... + b_K(xK)_i$$

Indeed, one way to assess the success of the regression is the closeness of these fitted Y values, namely  $\hat{Y}_1, \hat{Y}_2, \hat{Y}_3, ..., \hat{Y}_n$  to the actual observed Y values  $Y_1, Y_2, Y_3, ..., Y_n$ .

#### THIS IS LOOKING COMPUTATIONALLY HOPELESS.

Indeed it is. These calculations should only be done by computer. Even a careful, well-intentioned person is going to make arithmetic errors if attempting this by a non-computer method. You should also be aware that computer programs seem to compete in using the latest innovations. Many of these innovations are passing fads, so don't feel too bad about not being up-to-the-minute on the latest changes.

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The notation used here in the models is not universal. Here are some other possibilities.

Notation here	Other notation
$Y_i$	$y_i$
$x_i$	$X_i$
$\beta_0 + \beta_1 x_i$	$\alpha + \beta x_i$
$\epsilon_i$	$e_i$ or $r_i$
$(x1)_i, (x2)_i, (x3)_i,, (xK)_i$	$x_{i1}, x_{i2}, x_{i3}, \ldots, x_{iK}$
$b_j$	$\hat{\beta}_i$

In many regression problems, the data points differ dramatically in gross size.

EXAMPLE 1: In studying corporate accounting, the data base might involve firms ranging in size from 120 employees to 15,000 employees.

EXAMPLE 2: In studying international quality of life indices, the data base might involve countries ranging in population from 0.8 million to 1,000 millions.

In Example 1, some of the variables might be highly dependent on the firm sizes. For example, the firm with 120 employees probably has low values for gross sales, assets, profits, and corporate debt.

In Example 2, some of the variables might be highly dependent on country sizes. For example, the county with population 0.8 million would have low values for GNP, imports, exports, savings, telephones, newspaper circulation, and doctors.

Regressions performed with such gross size variables tend to have very large  $R^2$  values, but prove nothing. In Example 1, one would simply show that big firms have big profits. In Example 2, one would show that big countries have big GNPs. The explanation is excellent, but rather uninformative.

There are two common ways for dealing with the gross size issue: ratios and logarithms.

The ratio idea just puts the variables on a "per dollar" or "per person" basis.

For Example 1, suppose that you wanted to explain profits in terms of number of employees, sales, assets, corporate debt, and (numerically coded) bond rating. A regression of profits on the other variables would have a high  $R^2$  but still be quite uninformative. A more interesting regression would create the dependent variable profits/assets and use as the independent variables employees/assets, sales/assets, debt/assets. The regression model is

$$\frac{PROFIT_{i}}{ASSETS_{i}} = \beta_{0} + \beta_{1} \frac{EMPLOYEES_{i}}{ASSETS_{i}} + \beta_{2} \frac{SALES_{i}}{ASSETS_{i}} + \beta_{3} \frac{DEBT_{i}}{ASSETS_{i}} + \beta_{4} BOND_{i} + \varepsilon_{i}$$
(Model 1)

Observe that BOND, the bond rating, is not a "gross size" variable; there is no need to scale it by dividing by ASSETS.

In Example 1, the scaling might be described in terms of quantities per \$1,000,000 of ASSETS. It might also be reasonable to use SALES as the scaling variable, rather than ASSETS.

For Example 2, suppose that you wanted to explain number of doctors in terms of imports, exports, savings, telephones, newspaper circulation, and inflation rate. The populations give you the best scaling variable. The regression model is

$$\frac{DOCTORS_{i}}{POPN_{i}} = \beta_{0} + \beta_{1} \frac{IMPORTS_{i}}{POPN_{i}} + \beta_{2} \frac{EXPORTS_{i}}{POPN_{i}} + \beta_{3} \frac{SAVINGS_{i}}{POPN_{i}} + \beta_{4} \frac{PHONES_{i}}{POPN_{i}} + \beta_{5} \frac{PAPERS_{i}}{POPN_{i}} + \beta_{6} INFLATE_{i} + \varepsilon_{i}$$
(Model 2)

All the ratios used here could be described as "per capita" quantities. The inflation rate is not a "gross size" variable and need not be put on a per capita basis.

An alternate strategy is to take logarithms of all gross size variables. In Example 1, one might use the model

$$\log(PROFIT_{i}) = \gamma_{0} + \gamma_{1} \log(ASSETS_{i}) + \gamma_{2} \log(EMPLOYEES_{i}) + \gamma_{3} \log(SALES_{i})$$
$$+ \gamma_{4} \log(DEBT_{i}) + \gamma_{5} BOND_{i} + \varepsilon_{i}$$

Of course, the coefficients  $\gamma_0$  through  $\gamma_5$  are not simply related to  $\beta_0$  through  $\beta_4$  in the original form of the model. Unless the distribution of values of BOND is very unusual, one would not replace it with its logarithm.

Similarly, the logarithm version of model 2 is

$$\log(DOCTORS_i) = \gamma_0 + \gamma_1 \log(POPN_i) + \gamma_2 \log(IMPORTS_i) + \gamma_3 \log(EXPORTS_i)$$
$$+ \gamma_4 \log(SAVINGS_i) + \gamma_5 \log(PHONES_i) + \gamma_6 \log(PAPERS_i) + \gamma_7 INFLATE_i + \varepsilon_i$$

Since INFLATE is not a "gross size" variable, we are not immediately led to taking its logarithm. If this variable has other distributional defects, such as being highly skewed, then we might indeed want its logarithm.

Finally, it should be noted that one does not generally combine these methods. After all, since  $\log\left(\frac{A}{B}\right) = \log(A) - \log(B)$  the logarithm makes the ratio a moot issue.

Dividing logarithms, as in  $log(DOCTORS_i)/log(POPN_i)$  is not likely to be useful.

One always has the option of doing a "weighted" regression. One can use one of the variables as a weight in doing the regression. The company assets might be used for Example 1 and the populations used for Example 2. The problem with this approach is that the solution will depend overwhelmingly on the large firms (or large countries).

Data cleaning steps

We will describe the operations in terms of the computer program Minitab.

We will assume here that we are working with a spreadsheet. The columns of this spreadsheet will represent variables; each number in a column must be in the same units. The rows of the spreadsheet will represent data points.

As a preliminary step, check each column for basic integrity. Minitab distinguishes columns of two major varieties, ordinary data and text. (There are also minor varieties, including dates.) If a column is labeled C5-T, then Minitab has interpreted this column as text information.

It sometimes happens that a column which is supposed to be numeric ends up as text. What should you do in such a case?

Scan the column to check for odd characters, such as N/A, DK, ?, unk; some people use markers like this to indicate missing or uncertain values. The Minitab missing numeric data code is the asterisk \*, and this should be used to replace things like the above. The expression 2 1/2 was intended to represent 2.5 but Minitab can only interpret it as text; this repair is obvious.

If you edit a text column so that all information is interpretable as numeric, Minitab will not instantly recognize the change. Use **Manipulate**  $\Rightarrow$  **Change Data Type**  $\Rightarrow$  **Text to Numeric**. If you do this to a column that still has text information, the corresponding entries will end up as \*, the numeric missing data code.

It sometimes happens that a column given as numeric really represents a nominal categorical variable and you would prefer to use the names. For example, a column might have used 1, 2, 3, 4 to represent single, married, widowed, and divorced. You would prefer the names. Use **Manipulate**  $\Rightarrow$  **Code**  $\Rightarrow$  **Numeric to Text**. You will be presented with a conversion table which allows you to do this.

The command  $Stat \Rightarrow Basic Statistics \Rightarrow Display Descriptive Statistics will give you the minimum and maximum of each column. The minimum and maximum values should make sense; unbelievable numbers for the minimum or the maximum could well be data coding errors. This same command will give you the number of missing values, noted as <math>N^*$ . The count on missing values should make sense.

For many analyses you would prefer to deal with reasonably symmetric values. One of the cures for right-skewness is the taking of logarithms. Here are some general comments about this process:

Base *e* logarithms are usually preferred because of certain advantages in interpretation. It is still correct, however, to use base 10 logarithms.

Some variables are of the "gross size" variety. The minimum to maximum span runs over several orders of magnitudes. For example, in a data set on countries of the world, the variable POPULATION will run from 10<sup>5</sup> to 10<sup>9</sup> with many countries at the low end of the scale. This variable should be replaced by its logarithm. In a data set on the Fortune 500 companies, the variable REVENUES will run over several orders of magnitude with most companies toward the low end of the scale. This variable should be replaced by its logarithm.

The command  $Stat \Rightarrow Basic Statistics \Rightarrow Display Descriptive Statistics will allow you to compare the mean and the standard deviation. If a variable which is always (or nearly always) positive has a standard deviation about as large as the mean, or even larger, is certainly positively skewed.$ 

What should you do with data that are skewed but not necessarily of the "gross size" variety? This is a matter of judgment. Generally you prefer to keep variables in their original units. If most of the other variables are to be transformed by logarithms, then maybe you want to transform this one as well.

If the skewed variable is going to be the dependent variable in a regression, then you will almost certainly want to take its logarithm. (If you don't take the logarithms immediately, you may find expanding residuals on the residual versus fitted plot. Then you'll have take logarithms anyhow.)

If the variable to be transformed by logarithms as zero or negative values, then taking logarithms in Minitab will result in missing values! This is not usually what you want to do. Pick a value c so that all values of X + c are positive. Then consider  $\log(X + c)$ .

Logarithms will *not* cure left-skewed data. If X is such a variable and if M is a number larger than the biggest X, then you can consider  $\log(M-X)$ , provided you can make a sensible interpretation for this.

Logarithms should *not* be applied to binary variables. If a variable has only two values, then the logarithms will also have only two values.

Suppose that we regress Y on other variables, including J. The fitted model will be

$$\hat{Y} = b_0 + \ldots + b_J J + \ldots$$

The interpretation of  $b_J$  is this:

As J increases by 1, there is an associated increase in Y of  $b_J$ , while holding all other predictors fixed.

There's an important WARNING.

WARNING: This interpretation should note that  $b_J$  is the "effect" of J on Y after adjusting for the presence of all other variables. (In particular, regressing Y on J without any other predictors could produce a very different value of  $b_J$ .) Also, this interpretation carries the disclaimer "while holding all other predictors fixed." Realistically, it may not be possible to change the value of J while leaving the other predictors unchanged.

Now... suppose that Y is really the base-e logarithm of Z, meaning  $Y = \log Z$ . What's the link between J and Z? The fitted model is

$$\log \hat{Z} = b_0 + \dots b_J J + \dots$$

Here the interpretation of  $b_J$  is this:

As J increases by 1, there is an associated increase in  $\log Z$  of  $b_J$ . This means that  $\log Z$  changes to  $\log Z + b_J$ . By exponentiating, we find that  $e^{\log Z} = Z$  changes to  $e^{\log Z + b_J} = e^{\log Z} e^{b_J} = Z e^{b_J}$ . Using the approximation that  $e^t \approx 1 + t$  when t is near zero, we find that Z changes (approximately) to  $Z(1+b_J)$ . This is interpretable as a percent increase. We summarize thus: as J increases by 1, there is an associated proportional increase of  $b_J$  in Z.

If, for example,  $b_J = 0.03$ , then as J increases by 1, the associated increase in Z is 3%.

This next case is encountered only rarely.

Next suppose that Y is *not* the result of a transformation, but that  $J = \log R$  is the base-e logarithm of variable R. What's the link between R and Y? Let's talk about increasing J by 0.01. (The reason why we consider an increase of 0.01 rather than an increase of 1 will be mentioned below.) Certainly we can say this:

The fitted model is  $\hat{Y} = b_0 + ... + b_J \log R + ...$ 

As  $J = \log R$  increases by 0.01, there is an associated increase in Y of 0.01  $b_J$ . Saying that J increases by 0.01 is also saying that  $\log R$  increases to  $\log R + 0.01$ . By exponentiating, we find that  $e^{\log R} = R$  changes to  $e^{\log R + 0.01} = e^{\log R} e^{0.01} = R e^{0.01} \approx R (1+0.01)$ , which is a 1% increase in R.

Here's the conclusion: as R increases by 1%, there is an associated increase in Y of 0.01  $b_J$ .

If, for example,  $b_J = 25,400$ , then a 1% increase in R is associated with an approximate increase in Y of 254.

We used an increase of 0.01 (rather than 1) to exploit the approximation  $e^{0.01} \approx 1.01$ .

Finally, suppose that both Y and J are obtained by taking logs. That is  $Y = \log Z$  and  $J = \log R$ . What is the link between R and Z? Suppose we consider J increasing by 0.01; as in the previous note, this is approximately a 1% change in R.

As J increases by 0.01, there is an associated change from Y to Y + 0.01  $b_J$ . As  $Y = \log Z$ , we see that Z changes (approximately) to Z(1+0.01  $b_J$ ). Thus: as R increases by 1%, we find that there is an associated change in Z of 0.01  $b_J$ , interpreted as a percent.

If, for example,  $b_J = 1.26$ , then a 1% increase in R is associated with an approximate increase of 1.26% in Z.

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This document points out an interesting misunderstanding about multiple regression. There can be serious disagreement between

the regression coefficient  $b_H$  in the regression  $\hat{Y} = b_0 + b_G X_G + b_H H$  and

the regression coefficient  $b_H$  in the regression  $\hat{Y} = b_0 + b_H H$ 

While most people would not expect the values of  $b_H$  to be identical in these two regressions, it is somewhat shocking as to how far apart they can be.

Consider this very simple set of data with n = 20:

G	Н	Y
73	7.3	3096
87	-6.0	3519
83	-3.7	3383
78	2.5	3261
82	-2.2	3360
80	0.7	3334
83	-2.9	3388
86	-6.2	3481
75	5.1	3120
82	-1.3	3378

G	Н	Y
80	0.8	3326
82	-2.4	3365
77	2.9	3215
81	-1.5	3306
79	1.1	3266
78	1.9	3229
76	3.5	3193
80	0.5	3315
80	-0.3	3280
81	-0.6	3335

Here is the regression of Y on (G, H):

The regression equation is 
$$Y = -751 + 50.6 G + 20.5 H$$

Predictor	Coef	StDev	T	P
Constant	-751.2	515.9	-1.46	0.164
G	50.649	6.439	7.87	0.000
H	20.505	6.449	3.18	0.005

$$S = 13.63$$
  $R-Sq = 98.5\%$   $R-Sq(adj) = 98.3\%$ 

Analysis of Variance

Source	DF	SS	MS	F	P
Regression	2	209106	104553	562.64	0.000
Error	17	3159	186		
Total	19	212265			

This shows a highly significant regression. The F statistic is enormous, and the individual t statistics are positive and significant.

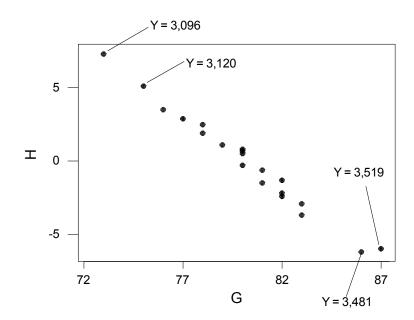
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Now, suppose that you regressed Y on H only. You'd get the following:

This regression is also highly significant. However, it now happens that the relationship with H is significantly negative.

How could this possibly happen? It turns out that these data were strung out in the (G, H) plane with a negative relationship. The coefficient of Y on G was somewhat larger than the coefficient on H, so that when we look at Y and H alone we see a negative relationship.

The picture below shows the locations of the points in the (G, H) plane. The values of Y are shown at some extreme points, suggesting why the apparent relationship between Y and H appears to be negative.



The quantity  $S_{yy} = \sum_{i=1}^{n} (y_i - \overline{y})^2$  measures variation in Y. Indeed we get  $s_y$  from this as  $s_y = \sqrt{\frac{S_{yy}}{n-1}}$ . We use the symbol  $\hat{y}_i$  to denote the fitted value for point i.

One can show that  $\sum_{i=1}^{n} (y_i - \bar{y})^2 = \sum_{i=1}^{n} (\hat{y}_i - \bar{y})^2 + \sum_{i=1}^{n} (y_i - \hat{y}_i)^2$ . These sums have the names  $SS_{\text{total}}$ ,  $SS_{\text{regression}}$ , and  $SS_{\text{error}}$ . They have other names or abbreviations. For instance

 $SS_{total}$  may be written as  $SS_{tot}$ .

 $SS_{regression}$  may be written as  $SS_{reg}$ ,  $SS_{fit}$ , or  $SS_{model}$ .

SS<sub>error</sub> may be written as SS<sub>err</sub>, SS<sub>residual</sub>, SS<sub>resid</sub>, or SS<sub>res</sub>.

The degrees of freedom accounting is this:

 $SS_{\text{total}}$  has n - 1 degrees of freedom

 $SS_{\text{regression}}$  has K degrees of freedom (K is the number of independent variables)

 $SS_{\text{error}}$  has n - 1 - K degrees of freedom

Here is how the quantities would be laid out in an analysis of variance table:

Source of Variation	Degrees of freedom	Sum of Squares	Mean Squares	F
Regression	K	$\sum_{i=1}^{n} (\hat{y}_i - \overline{y})^2$	$\frac{\sum_{i=1}^{n} (\hat{y}_i - \overline{y})^2}{K}$	$\frac{MS_{\text{Regression}}}{MS_{\text{Error}}}$
Error	n - 1 - K	$\sum_{i=1}^{n} \left( y_i - \hat{y}_i \right)^2$	$\frac{\sum_{i=1}^{n} (y_i - \hat{y}_i)^2}{n - 1 - K}$	
Total	n - 1	$\sum_{i=1}^{n} (y_i - \overline{y})^2$		

A measure of quality of the regression is the F statistic. Formally, this F statistic tests  $H_0: \beta_1 = 0, \beta_2 = 0, \beta_3 = 0, ..., \beta_K = 0$  [Note that  $\beta_0$  does not appear.] versus

 $H_1$ : at least one of  $\beta_1, \beta_2, \beta_3, ..., \beta_K$  is not zero

Note that  $\beta_0$  is not involved in this test.

Also, note that  $s_{\varepsilon} = \sqrt{MS_{\text{Error}}}$  is the estimate of  $\sigma_{\varepsilon}$ . This has many names:

standard error of estimate standard error of regression estimated noise standard deviation root mean square error (RMS error) root mean square residual (RMS residual)

The measure called  $R^2$  is computed as  $\frac{SS_{\text{Regression}}}{SS_{\text{Total}}}$ . This is often described as the "fraction of the variation in Y explained by the regression."

You can show, by the way, that

$$\frac{s_{\varepsilon}}{s_{v}} = \sqrt{\frac{n-1}{n-1-K} \left(1-R^{2}\right)}$$

The quantity  $R_{adj}^2 = 1 - \frac{n-1}{n-1-K} (1-R^2)$  is called the *adjusted R*-squared. This is supposed to adjust the value of  $R^2$  to account for both the sample size and the number of predictors. With a little simple arithmetic,

$$R_{adj}^2 = 1 - \left(\frac{s_{\varepsilon}}{s_{y}}\right)^2$$



This document considers the use of indicator variables, also called dummy variables, as predictors in multiple regression. Three situations will be covered.

EXAMPLE 1 gives a regression in which there are independent variables taking just two values. This is very easy.

EXAMPLE 2 gives a regression in which there is a discrete independent variable taking more than two values, but the values have a natural ordinal interpretation. This is also easy.

EXAMPLE 3 gives a regression with a discrete independent variable taking more than two values, and these values to not correspond to an ordering. This can get complicated.

#### EXAMPLE 1

Consider a regression in which the dependent variable SALARY is to be explained in terms of these predictors:

YEARS years on the job

SKILLS score on skills assessment (running from 0 to 40)

SUP 0 (not supervisor) or 1 (supervisor)

GENDER 0 (male) or 1 (female)

Suppose that the fitted regression turns out to be

Suppose that all the coefficients are statistically significant, meaning that the *p*-values listed with their *t* statistics are all 0.05 or less. We have these very simple interpretations:

The value associated with each year on the job is \$1,680 (holding all else fixed). The value associated with each additional point on the skills assessment is \$1,845 (holding all else fixed).

The value associated with being a supervisor is \$3,208 (holding all else fixed). The value associated with being female is -\$1,145 (holding all else fixed).

The variables SUP and GENDER have conveniently been coded 0 and 1, and this makes the interpretation of the coefficients very easy. Variables that have only 0 and 1 as values are called *indicator* variables or *dummy* variables.

If the scores for such a variable are two other numbers, say 5 and 10, you might wish to recode them.

These might also be described as categorical variables with two levels.

In general, we will not offer interpretations on estimated coefficients that are not statistically significant.

#### EXAMPLE 2

Consider a regression in which the dependent variable HOURS (television viewing hours per week) is to be explained in terms of predictors

```
INCOME (in thousands of dollars)

JOB (hours per week spent at work)

FAM (number of people living in the household)

STRESS (self-reported level of stress, coded as

1 = none, 2 = low, 3 = some, 4 = considerable, 5 = extreme)
```

The variable STRESS is clearly categorical with five levels, and we are concerned about how it should be handled. The important feature here is that STRESS is an *ordinal* categorical variable, meaning that the (1, 2, 3, 4, 5) responses reflect the exact ordering of stress. Accordingly, you need not take any extra action on this variable; you can use it in the regression exactly as is.

If the fitted regression equation is

$$HO\hat{U}RS = -62.0 - 1.1 INCOME - 0.1 JOB + 2.4 FAM - 0.2 STRESS$$

then the interpretation of the coefficient on STRESS, assuming that this coefficient is statistically significant, is that each additional level of STRESS is associated with 0.2 hour (12 minutes) less time watching television.

It seems natural to encode STRESS with consecutive integers. These are some subtleties:

- \* If you replaced the codes (1, 2, 3, 4, 5) by (-2, -1, 0, 1, 2), the regression would produce exactly the same estimated coefficient -0.2. This replacement would alter the intercept however.
- \* If you replaced the codes (1, 2, 3, 4, 5) by (10, 20, 30, 40, 50), the regression coefficient would be produced as -0.02.
- \* If you do not like the equal-size spaces between the codes, you might replace (1, 2, 3, 4, 5) by (-3, -1, 0, 1, 3). The coefficient would now change from -0.2, and you'd have to rerun the regression to see what it would be.

#### EXAMPLE 3

We will consider next a data set on home prices with n = 370.

Variable	Interpretation	Average	Standard deviation
PRICE	Home price in dollars	154,422	14,883
STYLE	Home style, coded as 1 = split-level, 2 = ranch, 3 = colonial, 4 = Tudor	2.41	0.98
SIZE	Indoor area in square feet	2,007.5	320.9
BEDROOM	Number of bedrooms	3.29	0.61

The number of bedrooms is a small integer, and we can use it in the regression with no modification. The average and standard deviation are useful summaries for BEDROOM, but we might also be interested in a simple tally. The following was obtained in Minitab from  $Stat \Rightarrow Tables \Rightarrow Tally$ .

BEDROOM	Count
2	21
3	230
4	109
5	10
N=	370

The variable STYLE is encoded as small integers, but the numbers function only as labels. Indeed, the information might have come to us as alphabetic names rather than these numbers. Note the inherent meaninglessness of the arithmetic

$$2-1$$
 = ranch – split-level =  $1 = 3-2$  = colonial – ranch

From **Stat**  $\Rightarrow$  **Tables**  $\Rightarrow$  **Tally** for the variable STYLE we get this:

STYLE	Count
1	85
2	97
3	141
4	47
N=	370

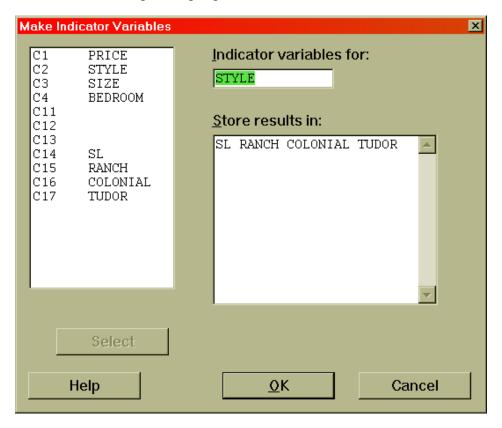
Since the numbers attached to STYLE do not mean anything, we cannot use this variable as presently structured.

By the way, if you uncritically ran the regression of PRICE on (STYLE, SIZE, BEDROOMS) you'd get the fitted equation

PRICE = 87443 + 5444 STYLE + 22.8 SIZE + 2436 BEDROOM and the coefficient on STYLE would be statistically significant. The interpretation would be that it's a \$5,444 step up from split-level to ranch, also a \$5,444 step up from ranch to Colonial, and a \$5,444 step up from ranch to Tudor. This is ridiculous.

If STYLE had only two values, we would be in the situation of EXAMPLE 1, and we could just use STYLE as an ordinary indicator (or dummy) variable. Here STYLE has four values, and we need a different method.

We will make a *set* of indicator variables for STYLE. In Minitab, do Calc  $\Rightarrow$  Make Indicator Variables. Set up the input panel to look like this:



In the box for **Store results in:** you must provide as many names as STYLE has values. Since STYLE has four values, we will give four names. For clerical sanity, we will match the names to the numeric values. That is, the 1-2-3-4 order will correspond directly to the names.

Minitab will assign the lowest number in STYLE to the first-named variable in the **Store results in:** list, the second-lowest number in STYLE to the second-named variable in the list, and so on. You need to be very careful. If you find this confusing, you might list a set of columns as C11-C14. After you see the results, you can assign names to these columns.

Minitab will create four indicator (dummy) variable columns. In the column for SL, the value 1 will appear for any house that was a split-level, and the value 0 will appear for all other houses. In the column for RANCH, the value 1 will appear for any house that was a ranch, and the value 0 will appear for all other houses.

In each row of the data sheet, SL + RANCH + COLONIAL + TUDOR will be exactly 1. This just notes that each house is one, and only one, of the four styles.

The command Calc ⇒ Make Indicator Variables can be applied to a column of alphabetic information. If STYLE contained alphabetic values, say

Split-Level Ranch Colonial Tudor then Minitab would assign

Colonial to the first-named variable in the **Store results in:** list Ranch to the second-named variable in the **Store results in:** list Split-Level to the third-named variable in the **Store results in:** list Tudor to the fourth-named variable in the **Store results in:** list This is done based on alphabetical ordering. Again, you need to be careful. You

still have the option of listing a set of columns as C11-C14. After you see the results, you can assign names to these columns.

It seems natural now to run the regression of PRICE on (SL, RANCH, COLONIAL, TUDOR, SIZE, BEDROOM). Note that STYLE is not included.

If you do that, you'll get this message at the top of the Minitab run:

- $^{\star}$  TUDOR is highly correlated with other X variables
- \* TUDOR has been removed from the equation

This message happens because SL + RANCH + COLONIAL + TUDOR = 1 for every line of the data set. This creates total collinearity with the regression intercept, and the regression arithmetic is impossible. Minitab deals with this by removing the last-named variable involved. In this instance, TUDOR was named last and was eliminated.

Minitab then goes on to produce a useful regression run:

```
The regression equation is
PRICE = 114696 + 21.8 SIZE + 2682 BEDROOM - 21054 SL - 12504 RANCH
            - 12639 COLONIAL
               SE Coef T P
114696 4160 27.57 0.000
21.832 1.993 10.96 0.000
2682 1006 2.66 0.008
-21054 1871 -11.26 0.000
-12504 1821 -6.86 0.000
-12639 1705 -7.41
              Coef SE Coef
114696 4160
Predictor
Constant
SIZE
BEDROOM
RANCH
COLONIAL
S = 9882
                 R-Sq = 56.5\% R-Sq(adj) = 55.9\%
Analysis of Variance
Source
                                  SS
                                                            F
Regression 5 46184185424 9236837085
                                                       94.58 0.000
Residual Error 364 35546964282 97656495
                  369 81731149706
Total
```

Parts of the output have been omitted.

The question now is the interpretation of the coefficients. For a split-level home, the indicators have values SL = 1, RANCH = 0, COLONIAL = 0. (Note that TUDOR has been omitted by Minitab). The fitted equation for a split-level home is then

```
PRICE = 114696 + 21.8 SIZE + 2682 BEDROOM - 21054 Split-Level
```

A ranch home has indicators SL = 0, RANCH = 1, COLONIAL = 0. This gives the fitted equation

```
PRICE = 114696 + 21.8 SIZE + 2682 BEDROOM - 12504 Ranch
```

Similarly, the fitted equation for colonial homes is

```
PRICE = 114696 + 21.8 SIZE + 2682 BEDROOM - 12639 Colonial
```

What about the Tudor homes? These have SL = 0, RANCH = 0, COLONIAL = 0, so that the fitted equation for these is

```
PRICE = 114696 + 21.8 SIZE + 2682 BEDROOM Tudor
```

The omitted indicator, here TUDOR, gives the base for interpreting the other estimated coefficients.

The suggestion is that a split-level home sells for 21,054 less than a Tudor home, holding all other variables fixed. A ranch sells for 12,504 less than a Tudor home, holding all other variables fixed. It follows that a ranch sells for 21,054 - 12,504 = 8,550 more than a split-level, holding all other variables fixed.

If we had asked Minitab for the regression of PRICE on (SL, RANCH, TUDOR, SIZE, BEDROOM), we would have produced the following fitted equation:

```
PRICE = 102057 + 21.8 SIZE + 2682 BEDROOM - 8415 SL + 135 RANCH + 12639 TUDOR
```

This time the indicator for colonial was used as the baseline, and we see that the Tudor homes sell for 12,639 more than the colonial homes, holding all else fixed. Perfectly consistent.

	Estimated coefficients			
Indicators used in the regression	SL	RANCH	COLONIAL	TUDOR
SL, RANCH, COLONIAL	-21,054	-12,504	-12,639	
SL, RANCH, TUDOR	-8,415	135		12,639
SL, COLONIAL, TUDOR	-8,550		-135	12,504
RANCH, COLONIAL, TUDOR		8 550	8 415	21.054

The following display indicates exactly what happens as we change the baseline.

In all parts of this table, the other variables (SIZE, BEDROOM) were used as well.

All four lines of this table represent equivalent fits. All produce the same  $R^2$ , the same F statistic, and the same  $s_{\varepsilon}$  (S in Minitab). Moreover, the estimated coefficients on SIZE and BEDROOM will be the same in all four lines, as will the corresponding t statistics.

If you are using a set of indicator variables, and if you go through a variable-selection process to remove variables, you must keep the indicator set intact. In the context of this problem, that means that any fitted model must use either

three out of the four indicators

or

none of the indicators

The indicators only make solid good sense when used together.

The regression of PRICE on (SIZE, BEDROOM, SL, RANCH, COLONIAL) which we saw above had significant *t* statistics on all independent variables. We would not be tempted to remove any of them. Moreover, a stepwise regression would select all the predictors.

The regression of PRICE on (SIZE, BEDROOM, SL, RANCH, TUDOR) produces this:

```
The regression equation is
PRICE = 102057 + 21.8 SIZE + 2682 BEDROOM - 8415 SL + 135 RANCH
        + 12639 TUDOR
                      SE Coef
Predictor
               Coef
                                       Т
                                                Ρ
                         3674
Constant
             102057
21.832
                                    27.78
                                            0.000
SIZE
                         1.993
                                    10.96
                                            0.000
              2682
                                            0.008
BEDROOM
                          1006
                                    2.66
              -8415
                                            0.000
                          1365
                                    -6.16
SL
RANCH
               135
                          1309
                                    0.10
                                            0.918
TUDOR
              12639
                           1705
                                     7.41
                                            0.000
```

This suggests that we might remove the indicator for RANCH. Indeed, stepwise regression selects all the variables except RANCH.

So what's the problem? If we removed RANCH, the other estimated coefficients would change, and we would no longer be able to assess correctly the differences between the home styles.

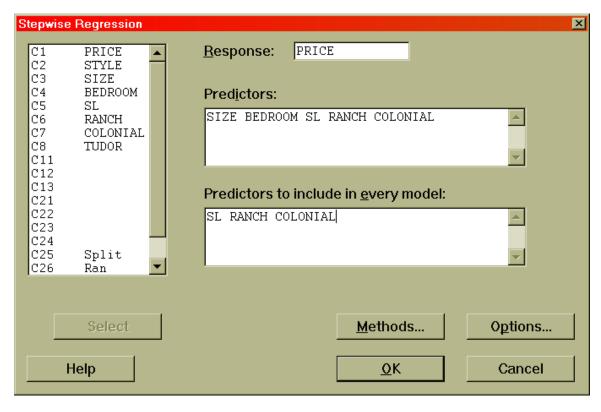
The advice, in generic form is this. If there are K indicators in a set, then a fitted model must use either

K-1 of the indicators (leave out any one)

or

none of the indicators.

Specifying a model that has none of the indicators is easy. If you use a variable selection technique like stepwise regression or best subsets regression, you need a way to force the indicator set to stay together. Here is how you set that up for stepwise regression:



Finally, we need an objective method to test whether an indicator variable set should be used at all. Let's consider the context of our model, namely

PRICE<sub>i</sub> = 
$$\beta_0$$
 +  $\beta_{SIZE}$  SIZE<sub>i</sub> +  $\beta_{BEDROOM}$  BEDROOM<sub>i</sub>  
+  $\beta_{SL}$  SL<sub>i</sub> +  $\beta_{RANCH}$  RANCH<sub>i</sub> +  $\beta_{COLONIAL}$  COLONIAL<sub>i</sub> +  $\epsilon_i$ 

The decision about whether or not to use the style indicators is really a test of the null hypothesis  $H_0$ :  $\beta_{SL} = 0$ ,  $\beta_{RANCH} = 0$ ,  $\beta_{COLONIAL} = 0$ .

There is a method for testing whether a *set* of coefficients is all zero. This method works for situations beyond what we are testing here. This requires the computation of this *F* statistic:

This is to be interpreted as an F statistic. We need to identify the two degrees of freedom numbers associated with F.

The numerator degrees of freedom is "Number of coefficients being investigated" in the calculation above.

The denominator degrees of freedom is the DF for residual in the regression on (SIZE, BEDROOM, SL, RANCH, COLONIAL).

The regression on (SIZE, BEDROOM, SL, RANCH, COLONIAL) had this analysis of variance table:

ance			
DF	SS	MS	F
5	46184185424	9236837085	94.58
364	35546964282	97656495	
369	81731149706		
	DF 5 364	DF SS	DF SS MS 5 46184185424 9236837085 364 35546964282 97656495

The regression sum of squares is 46,184,185,424. The residual mean square is 97,656,495. We note also that the degrees of freedom in the residual line is 364.

The regression on just (SIZE, BEDROOM) will have this analysis of variance table:

Analysis of Varia	ance			
Source	DF	SS	MS	F
P				
Regression	2	33675069487	16837534743	128.59
0.000				
Residual Error	367	48056080220	130942998	
Total	369	81731149706		

The regression sum of squares is 33,675,069,487.

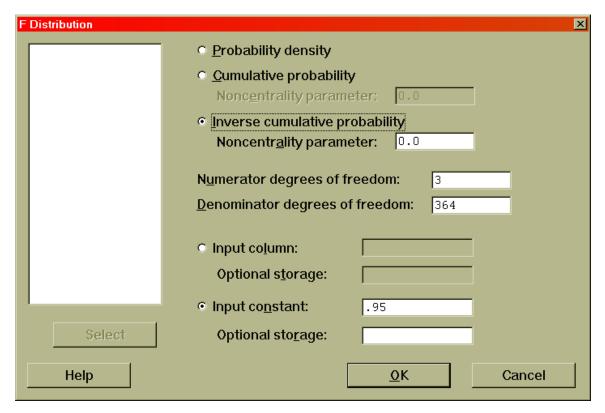
We'll note that three coefficients are under test. We now have enough information to assemble the test statistic:

$$\frac{\left\{46,184,185,424 - 33,675,069,487\right\} \div 3}{97,656,495} \approx 42.70$$

Minitab does not have a procedure for computing this number. The user needs to assemble it.

So what do we do with this number? The null hypothesis above should be rejected at the 0.05 level of significance if this exceeds  $F_{3,364}^{0.05}$ , the upper 5% point for the F distribution with (3, 364) degrees of freedom. It happens that  $F_{3,364}^{0.05} = 2.6294$ . Since our computed statistic, 42.70 exceeds 2.6294, we would reject the null hypothesis that all the coefficients of the style indicators are zero. It appears that the style indicators are useful as predictors of home price.

You can find this cutoff point for the F distribution from Minitab. Just do Calc  $\Rightarrow$  **Probability Distributions**  $\Rightarrow$  **F**, and then fill in the resulting panel as follows:



This particular F test had been defined through this statistic:

You will sometimes see this in the exactly equivalent form

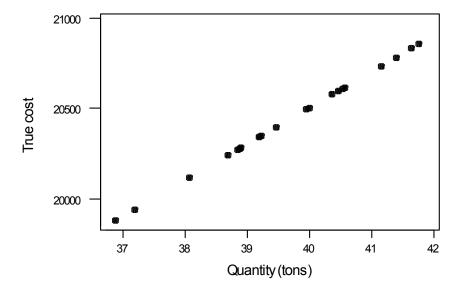
This equivalent form lays out the arithmetic as

$$\frac{\left\{48,056,080,220\ -\ 35,546,964,282\right\}\ \div\ 3}{97,656,495}\ \approx\ 42.70$$

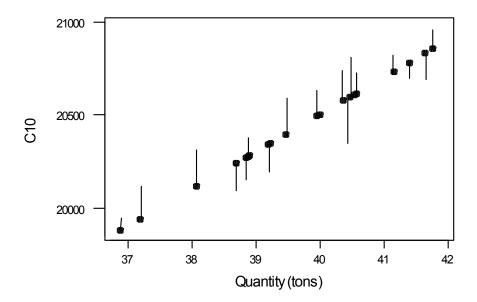
This produces exactly the same number, as it must.

There are many contexts in which regression analysis is used to estimate fixed and variable costs for complicated processes. The following data set involves the quantities produced and the costs for the production of a livestock food mix for each of 20 days. The quantities produced were measured in the obvious way, and the costs were calculated directly as labor costs + raw material costs + lighting + heating + equipment costs. The equipment costs were computed by amortizing purchase costs over the useful lifetimes, and the other costs are reasonably straightforward.

In fact, the actual fixed cost (per day) was \$12,500, and the variable cost was \$200/ton. Thus the exact relationship we see should be  $Cost = $12,500 + 200 \frac{\$}{ton} \times Quantity$ . Here is a picture of this exact relationship:



It happens, however, that there is statistical noise in assessing cost, and this noise has a standard deviation of \$100. Schematically, we can think of our original picture as being spread out with vertical noise:



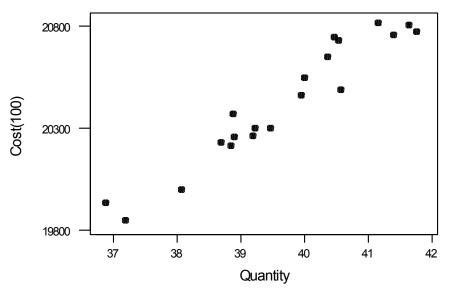
Here then are the data which we actually see:

Quantity	Cost	Quantity	Cost
41.66	20812.70	39.22	20302.30
40.54	20734.90	41.78	20776.70
38.90	20258.70	38.88	20373.00
38.69	20232.40	38.84	20213.70
40.58	20493.40	37.18	19848.70
40.48	20750.30	41.16	20818.90
36.88	19932.80	39.19	20265.10
39.47	20303.70	40.38	20654.50
41.41	20760.30	40.01	20553.00
38.07	20002.20	39.96	20463.10

The quantities are in tons, and the costs are in dollars.

Here is a scatterplot for the actual data:

## Costs in dollars to produce feed quantities in tons



(There is a noise standard deviation of \$100 in computing costs.)

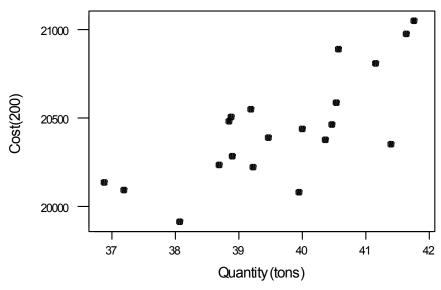
The footnote shows that in the process of assessing costs, there is noise with a standard deviation of \$100. In spite of this noise, the picture is fairly clean. The fitted regression line is  $\hat{Cost} = \$12,088 + 210 \frac{\$}{ton} \times \text{Quantity}$ . The value of  $R^2$  is 92.7%, so we know that this is a good regression. We would assess the daily fixed cost at \$12,088, and we would assess the variable cost at \$210/ton. Please bear in mind that this discussion hinges on knowing the exact fixed and variable costs and knowing about the \$100 noise standard deviation; in other words, this is a simulation in which we really know the facts. An analyst who sees only these data would not know the exact answer. Of course, the analyst would compute  $s_{\epsilon} = \$83.74$ , so that

Quantity	True value	Value estimated from data
Fixed cost	\$12,500	$b_0 = $12,088$
Variable cost	\$200/ton	$b_1 = \$210/\text{ton}$
Noise standard deviation	\$100	$s_{\varepsilon} = \$83.74$

All in all, this is not bad.

As an extension of this hypothetical exercise, we might ask how the data would behave with a \$200 standard deviation associated with assessing costs. Here is that scatterplot:

### Cost in dollars to produce feed quantities in tons



(There is a noise standard deviation of \$200 in computing costs.)

For this scatterplot, the fitted regression equation is  $\hat{Cost} = \$13,910 + 165 \frac{\$}{ton} \times \hat{Quantity}$ . Also for this regression we have  $R^2 = 55.4\%$ . Our estimates of fixed and variable costs are still statistically unbiased, but they are infected with more noise. Thus, our fixed cost estimate of \$13,910 and our variable cost estimate of  $165 \frac{\$}{ton}$  are not all that good. Of course, one can overcome the larger standard deviation in computing the cost by taking more data. For this problem, the analyst would see  $s_{\varepsilon} = \$210.10$ .

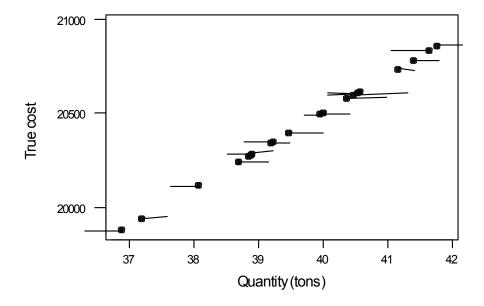
Quantity	True value	Value estimated from data
Fixed cost	\$12,500	$b_0 = \$13,910$
Variable cost	\$200/ton	$b_1 = $165/\text{ton}$
Noise standard deviation	\$200	$s_{\varepsilon} = \$210.10$

This is not nearly as good as the above, but this may be more typical.

It is important to note that noise in assessing cost, the vertical variable, still gives us a statistically valid procedure. The uncertainty can be overcome with a larger sample size.

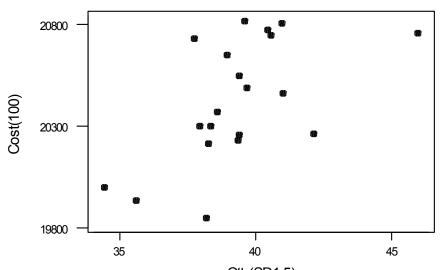
We will now make a distinction between noise in the vertical direction (noise in computing cost) and noise in the horizontal direction (noise in measuring quantity).

A more serious problem occurs when the horizontal variable, here quantity produced, is not measured exactly. It is certainly plausible that one might make such measuring errors when dealing with merchandise such as livestock feed. For these data, the set of 20 quantities has a standard deviation of 1.39 tons. This schematic illustrates the notion that our quantities, the horizontal variable, might not be measured precisely:



Here is a picture showing the hypothetical situation in which costs experienced a standard deviation of measurement of \$200 while the feed quantities had a standard deviation of measurement of 1.5 tons.

## Cost in dollars to produce feed quantities in tons



(There is a noise standard deviation of \$100 in computing costs) and quantities have been measured with a SD of 1.5 tons.)

For this picture the relationship is much less convincing. In fact, the fitted regression equation is  $\hat{\text{Cost}} = \$17,511 + 74.2\frac{\$}{\text{ton}} \times \text{Quantity}$ . Also, this has  $s_{\epsilon} = \$252.60$ . This has not helped:

Quantity	True value	Value estimated from data
Fixed cost	\$12,500	$b_0 = \$17,511$
Variable cost	\$200/ton	$b_1 = \$74.20/\text{ton}$
Noise standard deviation	\$200	$s_{\varepsilon} = $252.60$

The value of  $R^2$  here is 34.0%, which suggests that the fit is not good.

Clearly, we would like both cost and quantity to be assessed perfectly. However,

noise in measuring costs leaves our procedure valid (unbiased) but with imprecision that can be overcome with large sample sizes

noise in measuring quantities makes our procedure biased

The data do not generally provide clues as to the situation.

Here then is a summary of our situation.

Suppose that the relationship is

True cost = 
$$\beta_0 + \beta_1 \times$$
 True quantity

where  $\beta_0$  is the fixed cost and  $\beta_1$  is the variable cost

Suppose that we observe

$$Y = \text{True cost} + \varepsilon$$

where  $\epsilon$  represents the noise in measuring or assessing the cost, with standard deviation  $\sigma_\epsilon$ 

and

$$x = \text{True quantity} + \zeta$$

where  $\zeta$  represents the noise in measuring or assessing the quantity, with standard deviation  $\sigma_{\zeta}$ 

Let us also suppose that the True quantities themselves are drawn from a population with mean  $\mu_x$  and standard deviation  $\sigma_x$ .

You will do least squares to find the fitted line  $\hat{Y} = b_0 + b_1 x$ .

It happens that  $b_1$ , the sample version of the variable cost, estimates  $\beta_1 \frac{\sigma_x^2}{\sigma_x^2 + \sigma_r^2}$ .

Of course, if  $\sigma_{\zeta} = 0$  (no measuring error in the quantities), then  $b_1$  estimates  $\beta_1$ . It is important to observe that if  $\sigma_{\zeta} > 0$ , then  $b_1$  is biased closer to zero.

It happens that  $b_0$ , the sample version of the fixed cost, estimates

$$\beta_0 + \beta_1 \, \mu_x \, \frac{\sigma_\zeta^2}{\sigma_x^2 + \sigma_\zeta^2} \, .$$

If  $\sigma_{\zeta} = 0$ , then  $b_0$  correctly estimates the fixed cost  $\beta_0$ .

The impact in accounting problems is that we will tend to *underestimate* the variable cost and *overestimate* the fixed cost.

You can see that the critical ratio here is  $\frac{\sigma_{\zeta}^2}{\sigma_{x}^2}$ , the ratio of the variance of the noise in x relative to the variance of the population from which the x's are drawn.

In the real situation, you've got one set of data, you have no idea about the values of  $\beta_0$ ,  $\beta_1$ ,  $\sigma_x$ ,  $\sigma_{\zeta_0}$ , or  $\sigma_{\epsilon}$ . If you have a large value of  $R^2$ , say over 90%, then you can be pretty sure that  $b_1$  and  $b_0$  are useful as estimates of  $\beta_1$  and  $\beta_0$ . If the value of  $R^2$  is not large, you simply do not know whether to attribute this to a large  $\sigma_{\epsilon}$ , to a large  $\sigma_{\zeta}$ , or to both.

	Small $\sigma_{\zeta}/\sigma_{x}$ (quantity measured precisely relative to its background variation)	Large $\sigma_{\zeta}/\sigma_{x}$ (quantity measured imprecisely relative to its background variation)
Small $\sigma_{\epsilon}$ (cost measured precisely)	$b_0$ and $b_1$ nearly unbiased with their own standard deviations low; $R^2$ will be large	$b_1$ seriously biased downward and $b_0$ seriously biased upward; $R^2$ will not be large
Large $\sigma_{\epsilon}$ (cost measured imprecisely)	$b_0$ and $b_1$ nearly unbiased but their own standard deviations may be large; $R^2$ will not be large	$b_1$ seriously biased downward and $b_0$ seriously biased upward; $R^2$ will not be large

Do you have any recourse here?

If you know or suspect that  $\sigma_{\epsilon}$  will be large, meaning poor precision is assessing costs, you can simply recommend a larger sample size.

If you know or suspect that  $\sigma_{\zeta}$  will be large relative to  $\sigma_{x}$ , there are two possible actions:

By obtaining multiple readings of x for a single true quantity, it may be possible to estimate  $\sigma_{\zeta}$  and thus undo the bias. You will need to obtain the services of a serious statistical expert, and he or she should certainly be well paid.

You can spread out the x-values so as to enlarge  $\sigma_x$  (presumably without altering the value of  $\sigma_{\zeta}$ ). In the situation of our animal feed example, it may be procedurally impossible to do this.

### CROSS REF LIST (NOT FOR DISTRIBUTION)

## Original documents (not part of the formal handout)

introthoughts.doc
grossSize.doc
DataCleaning.doc
coefintr.doc
regpath.doc
anova.doc
indicator.doc (has a few things on variable selection)
errinx.doc