Correlation and Regression

references:

A&B Ch 5,8,9,10; Colton Chapter 6, M&M Chapters 2 (descriptive) and 9 (inference)

Similarities

- Both involve relationships between pair of numerical variables X & Y.
- Both: "predictability", "reduction in uncertainty"; "explanation".
- Both involve straight line relationships [can get fancier too].

Differences

- Correlation is **Symmetric**; Regression is **Directional**.
- Correlation involves sampling pairs of (X,Y) points; regression may involve choosing X values and sampling Y values to form the (X,Y) pairs or it may involve sampling pairs, but regards the X values as 'fixed'; x taken to be without error (not always so).
- Regression can be extended to relationships that are more complicated than just a straight line and can relate Y to multiple X variables.

Correlation [Pearson Product Moment Correlation Coefficient]

Loose Definition:

Degree to which, in observed (x,y) pairs, y value tends to be larger than average when x is larger (smaller) than average; extent to which larger than average x's are associated with larger (smaller) than average y's

Mathematically:

for (x,y) sample pairs,
$$r_{xy} = \frac{\{x_i - x\}\{y_i - y\}}{\sqrt{(\{x_i - \bar{x}\}^2)(\{y_i - \bar{y}\}^2)}}$$

"universe" of pairs: $\rho_{xy} = \frac{E\{(X - \mu_X)(Y - \mu_Y)\}}{\sqrt{E\{(X - \mu_X)^2\} E\{(Y - \mu_Y)^2\}}}$

<u>Positive Correlation:</u> Large X's with large Y's and small X's with small Y's <u>Negative Correlation:</u> Large X's with small Y's and small X's with large Y's <u>No Correlation:</u> Large X's with either large Y's or small Y's.



How r ranges from -1 (neg correln.) through 0 (zero correln.) through +1 (positive correln.) (r not tied to x or y scale).





 ρ^2 is a measure of how much the variance of Y is reduced by knowing what the value of X is (or vice versa) See article (by Chatillon) on "Balloon Rule" for visually estimating r.

 $Var(Y | X) = Var(Y) \cdot (1 - 2)$ $Var(X | Y) = Var(X) \cdot (1 - 2)$ $^{2} : "coefficient of determination"$

Large ² (i.e. close -1 or +1) -> close linear association of X and Y values; far less uncertain about value of one variable if told value of other. If X and Y scores are standardized to have mean=0 and unit SD=1 it can be seen that is like a "rate of exchange" ie the value of a standard deviation's worth of X in terms of PREDICTED standard deviation units of Y. If we know observation is Z_X SD's from μ_X , least squares prediction of observation's Z_Y value is given by **predicted** $Z_Y = \rho \cdot Z_X$ Notice the regression towards mean: is always less than 1 in absolute value, and so will make the predicted Z_Y closer to 0 (or equivalently make Y closer to to μ_Y) than the Z_X was to 0 (or X was to μ_X).

Inferences re ρ [based on sample of n (x,y) pairs]

Naturally, the observed r in any particular sample will not exactly match the ρ in the population (i.e. the coefficient one would get if one included <u>everybody</u>). The quantity r varies from one sample of n to another sample of n. i.e. r is subject to sampling fluctuations about ρ .

1. A common question asked of one's data is whether there is evidence of a <u>non-zero correlation</u> between 2 variables. To test this, one sets up a null hypothesis that ρ is zero and determines the probability, calculated under this null hypothesis that in fact $\rho = 0$, of obtaining an r more extreme than we observed. If the null is true, r would just be "randomly diffrent" from zero, with the amount of the random variation governed by n.

This discrepancy of r from zero is measured as $\frac{r \sqrt{n-2}}{\sqrt{1-r^2}}$ and should, if

the null hypothesis of $\rho = 0$ is true, follow t distribution with n-2 df. [Colton's table A5 gives the smallest r which would be considered evidence that ρ is not equal to 0. For example, if n=20, so that df = 18, an observed correlation of 0.44 or higher, or between -0.44 and -1 would be considered statistically significant at the P=0.05 level (2-sided). NB: the ttest assumes that the pairs are from a Bivariate Normal distribution. Also, it is valid only for testing $\rho = 0$, not for testing any other value of ρ .

Other common questions: given that r is based only on a sample, what <u>confidence interval</u> should I put around r so that there is a good chance (say 95%) that the interval will include the "true" coefficient ρ?

Or (answerable by the same technique): one observes a certain r_1 ; in another population, one observes a value r_2 . Is there evidence that the ρ 's in the 2 populations we are studying are unequal?

From our experience with the binomial statistic, which is limited to $\{0,n\}$ or $\{0,1\}$, it is no surprize that the r statistic, limited as it is to $\{-1,1\}$, also have a pattern of sampling variation that is <u>not be symmetric</u> unless ρ is right in the middle, i.e. unless $\rho = 0$. The following <u>transformation</u> of r will lead to a statistic which is approximately normal even if the ('s) in the population(s) we are studying is(are) quite distant from 0:

$$\frac{1}{2}$$
 ln { $\frac{1+r}{1-r}$ } [where ln is log to the base e or natural log].

It is known as Fisher's transformation of r; the observed r, transformed to this new scale, should be compared against a Gaussian distribution with

mean =
$$\frac{1}{2} ln \{ \frac{1+\rho}{1-\rho} \}$$
 and SD = $\sqrt{\frac{1}{n-3}}$.

<u>e.g. 2a</u>: **H**₀: $\rho = 0.5$; r=0.4 in sample of n=20. To test H₀, compute

$$\frac{\frac{1}{2} \ln \left\{ \frac{1+0.4}{1-0.4} \right\} - \frac{1}{2} \ln \left\{ \frac{1+0.5}{1-0.5} \right\}}{\sqrt{\frac{1}{n-3}}}$$
 and compare with

Gaussian (0,1) tables. Extreme values of the standardized Z are taken as evidence against H_0 . Often, the alternative hypothesis concerning ρ is 1-sided, of the form $\rho > some quantity$.

<u>e.g. 2b</u>: testing **H**₀: $\rho_1 = \rho_2$. $r_1 \& r_2$ in independent samples of $n_1 \& n_2$

Remembering that "variances add; SD's do not", compute the test statistic

$$\frac{\frac{1}{2} \ln \{\frac{1+r_1}{1-r_1}\} - \frac{1}{2} \ln \{\frac{1+r_2}{1-r_2}\} - [0]}{\sqrt{\frac{1}{n_1-3} + \frac{1}{n_2-3}}}.$$

and compare with Gaussian (0,1) tables.

e.g. 2c:
$$100(1-)$$
% **CI**(ρ) from r=0.4 in sample of n=20.

By solving the double inequality

$$-z_{1/2} = \frac{\frac{1}{2} \ln \left\{ \frac{1+r}{1-r} \right\} - \frac{1}{2} \ln \left\{ \frac{1+\rho}{1-\rho} \right\}}{\sqrt{\frac{1}{n-3}}} \qquad z_{1/2}$$

so that the middle term is ρ , we can construct a CI for $\rho \textbf{.}$

With a little re-arranging, it turns out to be ρ [lower, upper]

$$= \frac{1 + r - \{1 - r\} e^{\left[\pm 2 z^{/2} / \text{Sqrt[n-3]} \right]}}{1 + r + \{1 - r\} e^{\left[\pm 2 z^{/2} / \text{Sqrt[n-3]} \right]}}$$

<u>Worked e.g.</u> 95% CI(ρ) based on r=0.55 in sample of n=12.

With =0.05, so that $z_{1/2} = 1.96$, we have lower & upper bounds for ρ :

$$= \frac{1+0.55 - \{1-0.55\} e^{[\pm 2 \cdot 1.96 / \text{sqrt}[9]]}}{1+0.55 + \{1-0.55\} e^{[\pm 2 \cdot 1.96 / \text{sqrt}[9]]}}$$

$$= \frac{1.55 - 0.45 e^{[\pm 2 \cdot 1.96 / \text{sqrt}[9]]}}{1.55 + 0.45 e^{[\pm 2 \cdot 1.96 / \text{sqrt}[9]]}}$$

$$= \frac{1.55 - 0.45 e^{\pm 1.307}}{1.55 + 0.45 e^{\pm 1.307}} =$$

$$= \frac{1.55 - 0.45 \cdot 3.69}{1.55 + 0.45 \cdot 3.69} , \frac{1.55 - 0.45 / 3.69}{1.55 + 0.45 / 3.69} = -0.04 \text{ to}$$

0.84

This CI, which overlaps zero, agrees with the test of $\rho = 0$ described above. For if we evaluate $\frac{0.55\sqrt{12-2}}{\sqrt{1-0.55^2}}$, we get a value of 2.08,

which is not as extreme as the tabulated $t_{10,0.05(2-sided)}$ value of 2.23. <u>Note</u>: There will be some slight discrepancies between the t-test of $\rho = 0$ and the z-based CI's. The latter are only approximate. Note also that <u>both</u> assume we have data which have a bivariate Gaussian distribution.

A partial nomogram for 100(1–0.05)% i.e. 95% CI's for ρ is given on web page. It is based on Fisher's transformation of r. In addition to reading it vertically to get a CI for ρ (vertical axis) based on an observed r (horizontal axis), one can also use it to test whether an observed r is compatible with, or significantly different at the = 0.05 level, from some specific ρ value, ρ_0 say, on the vertical axis: simply read across from $\rho = \rho_0$ and see if the observed r falls within the horizontal range appropriate to the sample size involved. Note that this test of a <u>nonzero</u> ρ is not possible via the t-test. Books of statistical tables have fuller nomograms.

Spearman's (Non-parametric) Rank Correlation Coefficient

How Calculated:

• (i) replace the raw x's and y's by their ranks (1=smallest to n=largest)

(ii) calculate Pearson correlation on the ranks.

<u>Advantages</u>

• Easy to do manually (if ranking not a chore);

$$r_{\text{Spearman}} = 1 - \frac{6 \text{ d}^2}{n\{n^2 - 1\}}$$

- $\{ d = in ranks on X \& Y for each obsn. \}$
- Less sensitive to outliers $(x \rightarrow rank = variance fixed (for a given n).$ Extreme $\{x_i - \overline{x} \}$ or $\{y_i - \overline{y}\}$ can exert considerable influence on $r_{Pearson}$.
- Picks up on non-linear patterns e.g. the r_{Spearman} for the following data is 1, whereas the r_{Pearson}. is less.



Correlations -- obscured and artifactual



Regression

Uses			MAI	ES			FEMZ	ALES	
Curve fitting	Age.	Tot.	%-il	e; weig	ght,g	Tot.	%-il€	e; weigh	t,g
• Summarization ('model')	,		10.1	50.1	0.0.1		10.1	50.1	0.0.1
Description	wĸ	Ν.	lOth		90th	No.	lOth	50th	90th
Prediction									
• Explanation	25	100	651	810	<u>950</u>	73	<u>_604</u>	750	924
 Adjustment for 'confounding' variables 									
	30	257	<u>1 156</u>	1 530	<u>2_214</u>	216	<u>1_040</u>	1 485	<u>2_001</u>
Technical Meaning	31								
• [originally] simply a line of 'best fit' to data points	32								
	33								
• [nowadays] Regression line is the LINE that connects the CENTRES of	34	1 010	2 060	2 570	2 1 1 0	1 151	1 050	2 460	2 040
the distributions of Y's at each X value.	35	1 040	2_000	2 370	<u> </u>	1 454	1 950	<u>2 400</u>	5_040
	36								
• not necessarily a straight line; could be curved, as with growth charts	38								
	39								
 not necessarily µ_{Y X} 's used as CENTRES ; could use medians etc. 	40 6	8 102	3 020	3 570	4 160	67 149	2 900	3 430	4 000
	41								
• strictly speaking, haven't completed description unless we characterize	42 1	0 309	3 200	3 770	4 390	9 636	3 060	3 610	4 190
the variation around the centres of the Y distributions at each X									
		BIR	TH WEIG	HT (DIST	RIBUTIC	ON) MALE	S		
• inference not restricted to the distributions of Y's for which we make		BIR	TH WEIG	HT (MEC	DIAN) FEI	MÁLES			
some observations: it applies to distributions of Y's at all unobserved X				(EOth 0/:	la) for M/				
values in hetween	-	·	 Median Median 	(50th %) (50th %	ile) for FF	MALES			
values in between.			e meala	. (0011 /0			Ń		
Examples (with appropriate caveats)							` ` .		
• Dirth weight (V) in relation to gestational age (V)	4000a -					_			
 Blood prossure (Y) in relation to age (X) 	looog							· · ·	
 Cardiovascular mortality (V) in relation to water hardness (X) ? 					1)	
 Cancer incidence (V) in relation to some exposure (X)? 	3000a			į		o			
• Scholastic performance (Y) vis a vis amount of TV watched (X)	3000g -			}				/	
scholastic performance (1) his a his amount of 1 + mached (1)			k		Ŭ	/	/		

Caveat: No guarantee that simple straight line relationship will be adequate. Also, in some instances the relationship might change with the type of X and Y variables used to measure the two phenomena being studied; also the relationship may be more artifact than real - see later for inference.)

2000g

1000g

0

34

GESTATIONAL AGE (week)

36

38

Live singleton births, Canada 1986 Source: Arbuckle & Sherman CMAJ 140 157-161, 1989

40

 \cap

32

0

30

Simple Linear[†] Regression (one X) (straight line)

<u>Equation</u>

• $\mu_{Y|X} = + X$ or $\frac{\mu_{Y|X}}{X} = = \frac{"rise"}{"run"}$

In Practice:

one <u>rarely</u> sees an <u>exact</u> straight line relationship in health science applications;

- 1 While physicists are often able to examine the relationship between Y and X in a laboratory with <u>all other things being equal</u> (ie controlled or held constant) medical investigators largely are not. The universe of (X,Y) pairs is very large and any 'true' relationship is <u>disturbed</u> by <u>countless uncontrollable</u> (and sometimes <u>un-measurable factors</u>. In any particular <u>sample</u> of (X,Y) pairs these distortions will surely be operating.
- 2 The true relationship (even if we could measure it exactly) may not be a simple straight line.
- 3 The measuring instruments may be faulty or inexact (using 'instruments' in the broadest sense).

One always tries to have the investigation sufficiently controlled that the 'real' relationship won't be 'swamped' by factors 1 and 3 and that the background "noise" will be small enough so that alternative models (eg curvilinear relationships) can be distinguished from one another.

[†] Linear here means linear in the <u>parameters</u>. The equation $y = Bx^{C}$ can be made <u>linear in the parameters</u> by taking logs i.e. log[y] = log[B] + x log[C]; $y = a+b \cdot x+c \cdot x^{2}$ is already linear in the parameters a b and c. The following model cannot be made linear in the parameters :

proportion dying = $+ \frac{1}{1+\exp\{-\log[dose]\}}$

Fitting a straight line to data - Least Squares Method

The most common method is that of <u>Least Squares</u>. Note that least squares can be thought of as just a curve fitting method and doesn't have to be thought of in a statistical (or random variation or sampling variation) context. Other more statistically-oriented methods include the method of minimum Chi-Square (matching observed and expected counts according to measure of discrepancy) and the Method of Maximum likelihood (finding the parameters that made the data most likely). Each has a different criterion of "best-fit".

Least Squares Approach:

Consider a candidate slope (b) and intercept (a) and predict that the Y value accompanying any X=x is $\hat{y} = a + b \cdot x$. The <u>observed</u> y value will deviate from this "<u>predicted</u>" or "fitted" value by an amount d = y - \hat{y}

We wish to keep this deviation as <u>small</u> as <u>possible</u>, but we must try to strike a <u>balance</u> over all the data points. Again just like when calculating variances, it is easier to work with <u>squared</u> deviations¹:

$$f^2 = (y - \hat{y})^2$$

We <u>weight all deviations equally</u> (whether they be the ones in the middle or the extremes of the x range) using $d^2 = (y - \hat{y})^2$ to measure the overall (or average) discrepancy of the points from the line.

¹ there are also several theoretical advantages to least squares estimates over others based for example on least absolute deviations: - they are the most precise of all the possible estimates one could get by taking linear combinations of y's.

- From all the possible candidates for slope (b) and intercept (a), we choose the particular values a and b which make this sum of squares (sum of squared deviations of 'fitted' from 'observed' Y's) a minimum. ie we search for the a and b that give us the least squares fit.
- Fortunately, we don't have to use trial and error to arrive at the 'best' a and b. Instead, it can be shown by calculus or algebraically that the a and b which minimize d² are:

$$b = ^{a} = \frac{\{x_{i} - \bar{x}\}\{y_{i} - \bar{y}\}}{\{x_{i} - \bar{x}\}^{2}} = \frac{r_{xy} \cdot s_{y}}{s_{x}}$$
$$a = ^{a} = \bar{y} - b \bar{x}$$

[Note that a least-squares fit of the regression line of <u>X on Y</u> would give a <u>different</u> set of values for the slope and intercept: the slope of the line of x on y is $\frac{r_{Xy} \cdot s_X}{s_y}$. one needs to be careful when using a calculator or computer program to specify which is the explanatory variable (X) and which is the predicted variable (Y)].

Meaning of intercept parameter (a):

Unlike the <u>slope</u> parameter (which represents the increase/decrease in $\mu_{Y|X}$ for every unit increase in x), the intercept does not always have a 'natural' interpretation. It depends on where the x-values lie in relation to x=0, and may represent part of what is really the mean Y. For example, the regression line for fuel economy of cars (Y) in relation to their weight (x) might be

 $\mu_{Y|weight} = 60 \text{ mpg} - 0.01 \text{ weight in lbs} [0.01 \text{ mpg/lb}]$

but there are no cars weighing 0 lbs. It would be better to write the equation in relation to some 'central' value for weight e.g. 3500 lbs; then the same equation can be cast as

$$\mu_{Y|weight} - 25 = 0.01 \cdot (weight - 3500)$$

It is helpful for testing whether there is evidence of a non-zero slope to think of the simplest of all regression models, namely that which is a horizontal straight line

$$\mu_{\mathbf{Y}|\mathbf{X}} = + \mathbf{O} \cdot \mathbf{X} = \text{the constant}$$

This is a re-statement of the fact that the sum of squared deviances around a constant horizontal line at height 'a' is smallest when ' = the mean .

[We don't always use the mean as the best 'centre' of a set of numbers. Imagine waiting for one of several elevators with doors in a row along one wall; you do not know which one will arrive next, and so want to stand in the 'best' place no matter which one comes next. Where to stand depends on the criterion being optimized: if you want to minimize the <u>maximum</u> distance, stand in the middle between the one on the extreme left and the extreme right; if you wish to minimize the <u>average</u> distance , where do you stand?, If, for some reason, you want to minimize the <u>average squared</u> distance, where to stand? If the elevator doors are not equally spaced from each other, what then?]

The anatomy of a slope: some re-expressions

Consider the formula: slope = $b = \frac{\{x - xbar\}\{y - ybar\}}{\{x - xbar\}^2}$

Without loss of generality & for simplicity, assume ybar=0.

If we have 3 x's, 1 unit apart (e.g. $x_1=1$; $x_2=2$; $x_3=3$),

then... $x_1 - xbar = -1; x_2 - xbar = 0; x_3 - xbar = +1$

so slope = b =
$$\frac{\{-1\}y_1 + \{0\}y_2 + \{+1\}y_3}{\{-1\}^2 + \{0\}^2 + \{+1\}^2}$$

i.e. slope = $\frac{y_3 - y_1}{x_3 - x_1}$

Note that y_2 contributes to ybar and thus to an estimate of the average y (i.e. level) but not to the slope.

If 4 x's 1 unit apart (e.g. $x_1=1$; $x_2=2$; $x_3=3$; $x_4=4$), then,...

$$x_{1} - xbar = -1.5 \qquad x_{2} - xbar = -0.5$$

$$x_{3} - xbar = +0.5 \qquad x_{4} - xbar = +1.5$$
and so
slope =
$$b = \frac{\{-1.5\}y_{1} + \{-0.5\}y_{2} + \{+0.5\}y_{3} + \{+1.5\}y_{4}}{\{-1.5\}^{2} + \{-0.5\}^{2} + \{-0.5\}^{2} + \{-0.5\}^{2} + \{+1.5\}^{2}}$$
i.e. slope =
$$\frac{1.5\{y_{4} - y_{1}\}}{5} + \frac{0.5\{y_{3} - y_{2}\}}{5}$$
i.e. slope =
$$\frac{\frac{3}{2}\{y_{4} - y_{1}\}}{\frac{5}{3}\{x_{4} - x_{1}\}} + \frac{\frac{1}{2}\{y_{3} - y_{2}\}}{\frac{5}{1}\{x_{3} - x_{2}\}}$$
i.e. slope =
$$\frac{9}{10}\frac{\{y_{4} - y_{1}\}}{\{x_{4} - x_{1}\}} + \frac{1}{10}\frac{\{y_{3} - y_{2}\}}{\{x_{3} - x_{2}\}}$$

i.e. a weighted average of the slope from datapoints 1 and 4 and that from datapoints 2 and 3, with weights proportional to the squares of their distances on x axis $\{x_4 - x_1\}^2$ and $\{x_3 - x_2\}^2$

Another way to think of the slope:

Rewrite
$$b = \frac{\{x - xbar\}\{y - ybar\}}{\{x - xbar\}^2}$$
 as

$$b = \frac{\{x - xbar\} 2 \quad \frac{\{y - ybar\}}{\{x - xbar\}}}{\{x - xbar\} 2} = \frac{\text{weight} \quad \frac{\{y - ybar\}}{\{x - xbar\}}}{\text{weight}}$$

$$weight \quad \{x - xbar\}^2 \text{ for estimate} \quad \frac{\{y - ybar\}}{\{x - xbar\}} \text{ of slope}$$

Yet another way to think of the slope:

b is a weighted average of all the pairwise slopes
$$\frac{y_i - y_j}{x_i - x_j}$$

with weights proportional to $\{x_i - x_j\}^2$.

e.g. If 4 x's 1 unit apart

denote by $b_{1\&2}$ the slope obtained from $\{x_2, y_2\}$ & $\{x_1, y_1\}$, etc...

$$b = \frac{1.b_{1\&2} + 4.b_{1\&3} + 9.b_{1\&4} + 1.b_{1\&3} + 4.b_{2\&4} + 1.b_{3\&4}}{1+4+9+1+4+1} = 20$$

jh 6/94

Inferences regarding Simple Linear Regression

How reliable are

- (i) the (estimated) <u>slope</u>
- (ii) the (estimated) <u>intercept</u>
- (ii) the predicted mean Y at a given X
- (iv) the predicted y for a (future) individual with a given X

when they are based on data from a sample? i.e. how much would these estimated quantities change if they were based on a different random sample [with the same x values]?

We can use the concept of <u>sampling variation</u> to (i) <u>describe</u> the 'uncertainty' in our estimates via CONFIDENCE INTERVALS or (ii) carry out <u>TESTS of significance</u> on the parameters (slope, intercept, predicted mean).

We can describe the degree of reliability of (or, conversely, the degree of uncertainty in) an estimated quantity by the standard deviation of the possible estimates produced by different random samples of the same size from the same x's. We call this (obviously conceptual) S.D. the standard error of the estimated quantity (just like the standard error of the mean when estimating μ). helpful to think of slope as an average difference in means for 2 groups that are 1 x-unit apart.

The size of the standard error will depend on

- 1. how 'spread apart' the x's are
- 2. How good a fit the regression line really is (i.e. how small is the unexplained variation about the line)
- 3. How large the sample size, n, is.

Factors affecting reliability (in more detail)

1. <u>The spread of the X's:</u> The best way to get a reliable estimate of the slope is to take Y readings at X's that are quite a distance from each other. E.g. in estimating the "per year increase in BP over the 30-50 yr. age range", it would be better to take X=30,35, 40, 45,

50 than to take X = 38, 39, 49, 41, 42. Any individual fluctuations will 'throw off' the slope much less if the X's are far apart.



thick line: real (true) relation between average BP at age X and X : *thin lines*: possible apparent relationships because of individual variation when we study 1 individual at each of two ages (a) spaced closer together (b) spaced further apart.

Notes

Regression line refers to the relationship between the <u>average Y at a</u> <u>given X</u> to the <u>X</u>, and <u>not</u> to <u>individual Y's</u> vs X. Obviously of course if the individual Y's are close to the average Y, so much the better!

The above argument would suggest studying individuals at the extremes of the X values of interest. We do this if we are sure that the relationship is a linear one. If we are not sure, it is wiser -- if we have a choice in the matter -- to take a 3-point distribution.

There is a common misapprehension that a Gaussian distribution of X values is desirable for estimating a regression slope of Y on X. In fact, the 'inverted U' shape of the Gaussian is the least desirable!

Factors affecting reliability (continued)

 The (vertical) variation about the regression line: Again, consider BP and age, and <u>suppose</u> that indeed the average BP of <u>all</u> persons aged X + 1 is units higher than the average BP of all persons aged X, and that this linear relationship

average BP of persons aged $x = + \cdot X$ (average of Y's at a given $x = intercept + slope \cdot X$)

holds over the age span 30-50.

Obviously, <u>everybody</u> aged x=32 <u>won't</u> have <u>the exact same</u> BP, some will be above the average of 32 yr olds, some below. Likewise for the different ages x=30,...50. In other words, at any x there will be a distribution of y's about the average for age X. Obviously, how wide this distribution is about $+ \cdot X$ will have an effect on what slopes one could find in different samples (measure vertical spread around the line by)



thick line: real (true) relation between average BP at age X and X: *thin lines:* possible apparent relationships because of individual variation when we study 1 individual at each of two ages when the within-age distributions have (a) a narrow spread (b) a wider spread <u>NOTE:</u> For <u>unweighted</u> regression, should have roughly <u>same spread</u> <u>of Y's at each X.</u>

Factors affecting reliability (continued)

3. <u>Sample Size (n)</u> Larger n will make it more difficult for the types of extremes and misleading estimates caused by 1) poor X spread and 2) large variation in Y about $\mu_{Y|X}$, to occur. Clearly, it may be possible

to spread the x's out so as to maximize their variance (and thus reduce the n required) but it may not be possible to change the magnitude of the variation about $\mu_{\mathbf{Y}|\mathbf{X}}$ (unless there are other known factors

influencing BP). Thus the need for reasonably stable estimated \hat{y} [i.e.estimate of $\mu_{Y|X}$]

Standard Errors

SE(b) = SE([^]) =
$$\frac{1}{\sqrt{\{x_i - \bar{x}\}^2}}$$
;
SE(a) = SE([^]) = $\sqrt{\frac{1}{n} + \frac{\bar{x}^2}{\{x_i - \bar{x}\}^2}}$

(Note: there is a negative correlation between a and b).

<u>We don't usually know</u> so we estimate it from the data, using scatter of the y's from the fitted line i.e. SD of the residuals)

If examine the structure of SE(b), see that it reflects the 3 factors discussed above: (i) a large spread of the x's makes contribution of each observation to $\{x_i - \bar{x}\}^2$ large, and since this is in the denominator, it reduces the SE (ii) a small vertical scatter is reflected in a small and since this is in the numerator, it also reduces the SE of the estimated slope (iii) a large sample size means that $\{x_i - \bar{x}\}^2$ is larger, and like (i) this reduces the SE.

The formula, as written, tends to hide this last factor; note that $\{x_i - \bar{x}_i\}^2$ is what we use to compute the spread of a set of x's -- we simply divide it by n-1 to get a variance and then take the square root to get the sd. To make the point here, *simplify n-1 to n* and write

$$\{x_i - \bar{x}\}^2$$
 n•var(x), so that $\sqrt{\{x_i - \bar{x}\}^2}$ $\sqrt{n} \cdot sd(x)$

and the equation for the SE simplifies to (approx)

SE(b)
$$\frac{1}{\sqrt{n} \cdot sd(x)} = \frac{SD_{y|x} / SD_{x}}{\sqrt{n}}$$

with \sqrt{n} in its familiar place in the denominator of the SE (even in more complex SE's, this is where \sqrt{n} is usually found !)

The structure of SE(a): In addition to the factors mentioned above, all of which come in again in the expected way, there is the additional

factor of \bar{x}^2 ; since this is in the denominator, it increases the SE .

This is natural in that if the data, and thus \overline{x} , are far from x=0, then any imprecision in the estimate of the slope will project backwards to a large imprecision in the estimated intercept. Also, if one uses 'centered'

x's, so that $\overline{x} = 0$, the formula for the SE reduces to

$$SE(a) = \sqrt{\frac{1}{n}} = \frac{1}{\sqrt{n}}$$

and we recognize this as SE(\bar{y}) -- not surprisingly, since $\bar{y}\,$ is the 'intercept' for centered data.

CI's & Tests of Significance for $\stackrel{\wedge}{\alpha}$ and $\stackrel{\wedge}{\beta}$ are based on t-distribution (or Gaussian Z's if n large)

$$\alpha: \qquad \hat{\alpha} \pm \mathbf{t}_{n-2} \bullet \mathbf{SE}(\hat{\alpha})$$

H₀:
$$\alpha = \alpha_0$$
 $\mathbf{t}_{n-2} = \frac{\hat{\alpha} - \alpha_0}{\mathbf{SE}(\hat{\alpha})}$

 $\beta: \qquad \hat{\beta} \pm \mathbf{t}_{n-2} \cdot \mathbf{SE}(\hat{\beta})$

H₀:
$$\beta = \beta_0$$

t_{n-2} = $\frac{\hat{\beta} - \beta_0}{\mathbf{SE}(\hat{\beta})}$

1

Standard Error for Estimated $\mu_{Y|X}$ or 'average Y at X'

We estimate 'average Y at X' or $\mu_{Y|X}$ by $^{\wedge} + ^{\wedge} \cdot X$. Since the estimate is based on two estimated quantities, each of which is subject to sampling variation, it contains the uncertainty of both:

SE(estimated average Y at X) =
$$\sqrt{\frac{1}{n} + \frac{\{X - \bar{x}\}^2}{\{x_i - \bar{x}\}^2}}$$

Again, we must use an estimate $\hat{}$ of

First-time users of this formula suspect that it has a missing or an x instead of an xbar or something. There is no typographical error, and indeed if one examines it closely, it makes sense. X refers to the x-value at which one is estimating the mean -- it has nothing to do with the actual x's in the study which generated the estimated coefficients, except that the closer X is to the center of the data, the smaller the quantity $\{X - \bar{x}\}$ and thus the quantity $\{X - \bar{x}\}^2$, and thus the SE, will be. Indeed, if we estimate the **average Y** right at $X = \bar{x}$, the estimate is simply \bar{y} (since the fitted line goes through $[\bar{x}, \bar{y}]$) and its SE will be

$$\sqrt{\frac{1}{n} + \frac{\{\overline{x} - \overline{x}\}^2}{\{x_i - \overline{x}\}^2}} \quad \text{or} \quad \sqrt{\frac{1}{n}} = \frac{1}{\sqrt{n}} = SE(\overline{y}).$$

Confidence Interval for individual Y at X

A certain percentage P% of individuals are within tp • of the mean $\mu_{\mathbf{Y}|\mathbf{X}} = + \mathbf{\cdot} \mathbf{X}$, where tp is a multiple, depending on P, from the t or, if n is large, the Z table. However, we are not quite certain where exactly the mean $+ \mathbf{\cdot} \mathbf{X}$ is -- the best we can do is estimate, with a certain P% confidence, that it is within tp• SE($^{\wedge} + ^{\wedge} \mathbf{\cdot} \mathbf{X}$) of the point estimate $^{\wedge} + ^{\wedge} \cdot \mathbf{X}$. The uncertainty concerning the mean and the natural variation of individuals around the mean -- wherever it is -- combine in the expression for the estimated P% range of individual variation, which is as follows:

$$^{\wedge} + ^{\wedge} \cdot \mathbf{X} \pm \mathbf{t} \cdot \mathbf{\nabla} \sqrt{1 + \frac{1}{n} + \frac{\{\mathbf{X} - \overline{\mathbf{x}}\}^2}{\{\mathbf{x}_i - \overline{\mathbf{x}}\}^2}}$$

Both the CI for the estimated mean and the CI for individuals (ie the estimated percentiles of the distribution) are bow-shaped when drawn as

a function of **X**. They are narrowest at $\mathbf{X} = \overline{\mathbf{x}}$, and fan out from there. One needs to be careful not to confuse the much narrower CI for the mean with the much wider CI for individuals. If one can see the raw data, it is usually obvious which is which -- the CI for individuals is almost as wide as the raw data themselves.

cf. data on sleeping through the night; alcohol levels and eye speed.

Community Fluoride and Cavities Analysis via MYSTAT

Descriptive statistics

	DMFTX100 (Y)	FLPPM (X)	XY
N OF CASES	21	21	21
MEAN	537.381	0.700	228.038
VARIANCE	59236.648	0.556	30254.578
STANDARD DEV	243.386	0.746	173.938
SUM	11285.000	14.700	4788.800

Regression analysis

DEP	VAR:DMFTX100	N:21	MULT	R:	0.857	SOUARED	MULT.	R:	0.	734

ADJUSTED R²: 0.721 STANDARD ERROR OF ESTIMATE: 128.667

VARIABLE	COEFFICIENT	STD ERROR	STD COEF	Т	P(2 TAIL)
CONSTANT	733.198	38.959	0.000	18.820	0.000
FLPPM	-279.739	38.585	-0.857	-7.250	0.000

ANALYSIS OF VARIANCE

SOURCE	SUM-OF-SQUARES	DF	MEAN-SQUARE	F-RATIO	P
REGRESSION	870184.756	1	870184.756	52.563	0.000
RESIDUAL	314548.196	19	16555.168		



<u>Analysis via SAS</u>

DATA A; INPUT CITY 5-6 DMFTX100 10-13 FLPPM 18-201; CARDS;

PROC PRINT;	;			
	OBS	CITY	DMFTX100) FLPPM
	1	1	236	1.9
	2	2	246	2.6
	•	•		
	•	•	• • •	• • •
	20	20	823	0.1
	21	21	1037	0.1

PROC MEANS; VAR DMFTX100 FLPPM;

VARIABLE	N	MEA	N	STD	MIN.	MAX	STD	ERROR
				DEVN	VALUE	VALUE	OF	MEAN
DMFTX100	21	537.38	243.	3858	236.00	1037.0	53	.111
FLPPM	21	0.70	0.	7456	0.00	2.6	0	.162

PROC GLM; MODEL DMFTX100 = FLPPM;

GENERAL LIN	IEAR MODELS	PROCEDURE -	DEP. VARIAB	LE: DMFTX100
SOURCE	DF	SUM OF SQ.	MEAN SQUAR	<u>e fvalue</u>
MODEL	1	870184.75	870184.75	52.56
ERROR	19	314548.19	16555.16	PR > F
CORRECTED	TOTAL 20	1184732.95		0.0001
R-SQUARE	C.V.	ROO	I MSE DM	FTX100 MEAN
0.734499	23.9433	128.666	88857 5	37.38095238
SOURCE	DF	TYPE I	II SS F VAL	<u>ue pr > f</u>
FLPPM	1	870184.756	29497 52.	56 0.0001
		T FOR H0:	PR > T	STD ERROR OF
PARAMETER	ESTIMATE	PARAMETER	=0	ESTIMATE
INTERCEPT	733.1983	18.82	0.0001	38.959
FLPPM	-279.7392	7.25	0.0001	38.584