Week 6: Maximum Likelihood Estimation

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Outline

- An alternative way of estimating parameters: Maximum likelihood estimation (MLE)
- Simple examples: Bernoulli and Normal with no covariates
- Adding explanatory variables
- Variance estimation
- Intuition about the linear model using MLE
- Likelihood ratio tests, AIC, BIC to compare models
- Logit and probit with a latent variable formulation
- Predictions in logit and probit models

- Suppose that we know that the following ten numbers were simulated using a Bernoulli distribution: 0 0 0 1 1 1 0 1 1 1
- We can denote them by $y_1, y_2, ..., y_{10}$. So $y_1 = 0$ and $y_{10} = 1$
- Recall that the pdf of a Bernoulli random variable is $f(y; p) = p^{y}(1-p)^{1-y}$, where $y \in \{0, 1\}$
- The probability of 1 is p while the probability of 0 is (1 p)
- We want to figure out what is the p that was used to simulate the ten numbers. That's what we do not know
- What we do know is 1) they come from a Bernoulli distribution and 2) they are independent from each other, meaning that knowing one of them doesn't tell us anything about the other(s)

- Since we know the pdf that generated the numbers is Bernoulli, we know that the probability of the first number is $p^{y_1}(1-p)^{1-y_1}$
- The probability of the second is $p^{y_2}(1-p)^{1-y_2}$ and so on...
- We could replace the y_i with the actual numbers. For example, the first one is $y_1 = 0$ so the probability is just (1 p). I'll keep the symbols because we are going to make the problem more general in a bit
- What we do NOT know is the value of the parameter p
- Since we know that they are independent we could also write down the probability of observing all 10 numbers. That is, their joint probability
- Since they are independent their joint distribution is the multiplication of the 10 pdfs. Recall: $p(A \cap B) = P(A)P(B)$ if A and B are independent

- The joint probability is $[p^{y_1}(1-p)^{1-y_1}] \times [p^{y_2}(1-p)^{1-y_2}] \times \cdots \times [p^{y_10}(1-p)^{1-y_0}]$
- To make notation easier we'll use the product symbol \prod . For example, $\prod_{i=1}^{2} x_i = x_1 * x_2$
- So we can write the *joint probability* or the likelihood (L) of seeing those 10 numbers as:

$$L(
ho) = \prod_{i=1}^{10}
ho^{y_i} (1-
ho)^{1-y_i}$$

 Again, that is the probability that we observed the 10 numbers, although we do not know p yet

- Here comes the key insight. We can turn this problem into a maximization problem. What is the number p that maximizes the likelihood function L(p) given the observed data?
- That number will be the estimated parameter p that we'll call \hat{p}
- Yet another way: we want to find the \hat{p} that makes the joint likelihood of seeing those numbers as high as possible
- That's of course a calculus problem. We can take the derivative of L(p) with respect to p and set it to zero to find the optimal \hat{p}
- Of course, the second step is to verify that it's a maximum and not a minimum (take second derivative) and also verify that is unique, etc. We will skip those steps to keep this problem simple

- Taking that derivative is complicated because we would need to use the chain rule several times. A lot easier to make it a sum by taking the log; the log function is a **monotonic transformation**, it won't change the optimal \hat{p} value
- We will use several properties of the log, in particular: $log(x^ay^b) = log(x^a) + log(y^b) = a * log(x) + b * log(y)$
- We are going to generalize the problem by assuming we have *n* numbers instead of 10. It doesn't make it more difficult. We have: $lnL(p) = \sum_{i=1}^{n} y_i ln(p) + \sum_{i=1}^{n} (1 - y_i) ln(1 - p)$
- Which simplifies to: $lnL(p) = n\bar{y}ln(p) + (n n\bar{y})ln(1 p)$
- $(\bar{y} = \frac{\sum_{i=1}^{n} y_i}{n})$ This looks a lot easier; all we have to do is take $\frac{dln(p)}{dp}$, set it to zero, and **solve for p**

$$rac{dln(p)}{dp} = rac{nar{y}}{p} - rac{(n-nar{y})}{(1-p)} = 0$$

• After solving, we'll find that $\hat{p}(y_i) = \bar{y} = \sum_{i=1}^{n} \frac{y_i}{n}$

- So that's the MLE estimator of p. This is saying more or less the obvious: our best guess for the p that generated the data is the proportion of 1s, in this case p = 0.6, which is the same as $E[Y_i]$
- We would need to verify that our estimator satisfies the three basic properties of an estimator: bias, efficiency, and consistency (this will be in your exam)
- Note that we can plug in the optimal \hat{p} back into the ln likelihood function:
- $lnL(\hat{p}) = n\bar{y}ln(\hat{p}) + (n n\bar{y})ln(1 \hat{p}) = a$, where a will be a number that represents the **highest likelihood we can achieve** (we chose \hat{p} that way)
- **Drum roll**: We just discovered logistic regression

Example

• Simulated 100 Bernoulli rvs with p = 0.4

set obs 100 gen bernie = uniform()<0.4 sum bernie Variable | Obs Mean Std. Dev. Min Max ----bernie | 100 .46 .5009083 0 1 * We just showed that p hat is 0.46 * Let's get the highest value of the ln likelihood * Plug in p hat and the other values di 100*0.46*1n(0.46) + (100-100*0.46)*1n(1-0.46)-68,994376

 Again, we just did logistic regression "by hand." A logistic model with only a constant (no covariates), also known as the null model

Replicate using the logit command

We have used logit before

logit bernie Iteration 0: Iteration 1:	log likelih log likelih	ood = -68.994 ood = -68.994					
Logistic regre	ssion			Number	of obs	-	100
				LR chi2	(0)	=	-0.00
				Prob >	chi2	=	
Log likelihood	l = -68.99437	6		Pseudo R2 =		=	-0.0000
bernie	Coef.	Std. Err.				Conf.	Interval]
_cons	1603427	.2006431	-0.80	0.424	553	5959	.2329106
di 1/(1+exp(. .45999999	1603427))						

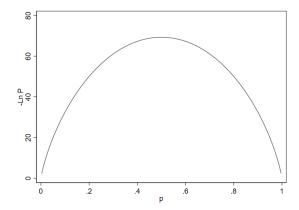
- Note that Log likelihood = -68.994376 matches what we found "by hand;" the coefficient is in the log-odds scale, but we can transform it
- This is a model with no explanatory variables. We can easily make the parameter p be a linear function of predictors

Some things to note

- Statistical software maximizes the log likelihood numerically. Also the log likelihood because of numerical precision. The probability of each y_i is small, so p(y₁) × p(y₂) × ... p(y_n) could be a very small number
- The algorithm is given a starting value for some parameters (often using using the null model)
- Each iteration "improves" the maximization
- The second derivatives are also computed because the second derivatives are used to compute SEs
- In many cases, we need to be mindful of the difference between the scale of estimation and the scale of interest
- Logit models report coefficients in the log-odds scale
- Not the only way of deriving logit models. You could also assume a Binomial pdf. The Bernoulli is a special case of the Binomial when the number of trials is 1. We will see a variable approach, too

Plot the - ln(L) function with respect to p

```
twoway function y = -(100^*x^*ln(x) + (100-100^*x)^*ln(1-x)), range(0 1) /// xtitle("p") ytitle("-Ln P") saving(1100.gph, replace)
```



What about the precision (standard error) of the estimate?

- There is some intuition in the plot above. The precision of the estimate \hat{p} can be measured by the curvature of the $lnL(\theta)$ function around its peak
- \blacksquare A flatter curve has more uncertainty. In logistic models, the variance is function of the p
- The Fisher information function, $I(\theta)$ formalizes that intuition: $I(\theta) = -E[\frac{\partial^2 lnL(\theta)}{\partial^2 \theta}]$
- It turns out that we can calculate $var(\theta)$ using the inverse of $I(\theta)$
- For the Bernoulli, $I(\hat{p}) = \frac{n}{[\hat{p}(1-\hat{p})]}$ (evaluated at \hat{p})
- The variance is $1/I(\hat{p}) = \frac{\hat{p}(1-\hat{p})}{n}$
- Note something. Once we know p̂ we also know its variance. The Normal distribution is unique in that the variance can change independently of the mean

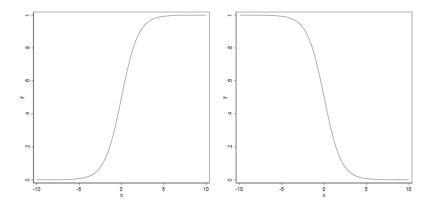
Logit models with covariates

- In most applications we want to estimate the effect of covariates on the probability p
- So we could just make p a function of covariates: $p = f(x_1, x_2, ..., x_p)$
- The easiest one is to make them a linear function like $p = \beta_0 + \beta_1 x_1 + \dots + \beta_p x_p$
- But that could be a problem since we need to guarantee that p will be bounded between 0 and 1 since it's a probability
- Enters the logistic or logit function: $\frac{1}{1+e^{-(\beta_0+\beta_1x_1+\cdots+\beta_px_p)}} = \frac{e^{(\beta_0+\beta_1x_1+\cdots+\beta_px_p)}}{1+e^{(\beta_0+\beta_1x_1+\cdots+\beta_px_p)}}$
- We have now changed the maximization problem. We do not maximize the likelihood with respect to p. Now we do it with respect to the parameters β_j
- The math is more difficult. We need matrix algebra. The second derivative becomes a matrix too (the Jacobian). The logic is the same, though

Logistic response function

 \blacksquare If we constrain the response to be between 0 and 1, it can't be linear with respect to X

```
twoway function y=\exp(x) / (1+\exp(x)), range(-10 10) saving(11.gph, replace)
twoway function y=\exp(-x) / (1+\exp(-x)), range(-10 10) saving(12.gph, replace)
graph combine 11.gph 12.gph, xsize(20) ysize(10)
graph export lboth.png, replace
```



Logistic or logit model

- Notice a couple of things. The effect of x on π is not linear; the effect depends on the value of x
- But we can make the function linear using the so-called logit transformation $ln(\frac{\pi}{1-\pi}) = x$
- I made you go the other way in one homework. If you solve for π you get to the logistic response function
- More general, the model is:

•
$$ln(\frac{p}{1-p}) = \beta_0 + \beta_1 X_1 + \dots + \beta_p X_p$$
, which transformed is
 $p = \frac{e^{\beta_0 + \beta_1 X_1 + \dots + \beta_p X_p}}{1 + e^{\beta_0 + \beta_1 X_1 + \dots + \beta_p X_p}}$

• Can also be written as: $p = \frac{1}{1 + e^{-(\beta_0 + \beta_1 X_1 + \dots + \beta_p X_p)}}$

What about if we do the same but now we have numbers like

90.46561 105.1319 117.5445 102.7179 102.7788 107.6234 94.87266 95.48918 75.63886 87.40594 ...

- I tell you that they were simulated from a normal distribution with parameters μ and σ^2 . The numbers are independent. Your job is to come up with the best guess for the two parameters
- Same problem as with the Bernoulli example. We can solve it in exactly the same way

As before, we know the pdf of a Normal random variable and because the observations are independent we can multiply the densities:

$$L(\mu, \sigma^2) = \prod_{i=1}^n \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(\frac{-(y_i - \mu)^2}{2\sigma^2}\right)$$

Remember the rules of exponents, in particular e^ae^b = e^{a+b}. We can write the likelihood as:

$$L(\mu, \sigma^2) = (\frac{1}{\sqrt{2\pi\sigma^2}})^n exp(-\frac{1}{2\sigma^2} \sum_{i=1}^n (y_i - \mu)^2)$$

- Alert: Perhaps you are wondering, why are we using the pdf of the normal if we know that the probability of one number is zero? Because we can think of the pdf as giving us the probability of $y_i + d$ when $d \rightarrow 0$
- If you recall your stats 101 classes, with a continuous rv the probability of a number is zero

After taking the In, we have:

$$lnL(\mu,\sigma^{2}) = -\frac{n}{2}ln(2\pi\sigma^{2}) - \frac{1}{2\sigma^{2}}\sum_{i=1}^{n}(y_{i}-\mu)^{2}$$

• All we have left is to take the derivative with respect to our two unknowns, μ and σ^2 and set them to zero. Let's start with μ :

$$\frac{\partial \ln(L(\mu,\sigma^2))}{\partial \mu} = 2\frac{1}{2\sigma^2}\sum_{i=1}^n (y_i - \mu) = 0$$

The above expression reduces to (I added the `to emphasize that's the optimal):

 $\sum_{i=1}^n (y_i - \hat{\mu}) = 0$

- **Does it look familiar?** Replace $\hat{\mu}$ with \hat{y}_i . That's exactly the same as the first order condition you saw when minimizing the sum of squares (that also the moment condition in Generalized Method of Moments, GMM)
- Solving, we find that $\hat{\mu} = \frac{\sum_{i=1}^{n} y_i}{n} = \bar{y}$. In other words, our best guess is just the mean of the numbers. $E[Y_i]$ again

- \blacksquare We can also figure out the variance by taking the derivative with respect to σ^2
- We will find that $\hat{\sigma}^2 = \frac{\sum_{i=1}^{n} (y_i \hat{\mu})}{n}$
- That formula happens to be wrong. That's a biased estimate of the variance. We need to divide by (n − 1) instead
- (What is the definition of bias?)
- This is not unusual in MLE. The MLE estimate of the variance is often biased but it is easy to correct them (and less of a problem with large samples)

Normal example Stata

- We just figured out that the best guess is to calculate the sample mean and sample variance
- We can easily verify in Stata

clear set seed 1234567 set obs 100 gen ynorm = rnor sum ynorm		0			
Variable	Obs	Mean			Max
ynorm	100		10.03931	74.16368	123.5079
reg ynorm < output omit	:ted>				
ynorm	Coef.		t P>		Conf. Interval]
_cons	98.52294			.000 96.5	3092 100.515

Drum roll 2: We just did linear regression "by hand" (although without covariates)

Linear regression: adding covariates

- What about if I told you that the number I generated is a linear function of one variable, say, x₁? In other words, I'm saying that the mean of the normal distribution is µ = β₀ + β₁x₁
- Now we want to find the parameters β₀, β₁, σ² that maximize the likelihood function. Once we know the optimal β̂₀, β̂₁ we find the optimal μ̂
- The likelihood function is now:

$$L(\beta_0, \beta_1, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} exp(-\frac{1}{2\sigma^2} \sum_{i=1}^n (y_i - \beta_0 - \beta_1 x_{1i})^2)$$

The In likelihood is:

$$lnL(\mu,\sigma^{2}) = -\frac{n}{2}ln(2\pi\sigma^{2}) - \frac{1}{2\sigma^{2}}\sum_{i=1}^{n}(y_{i} - \beta_{0} - \beta_{1}x_{1i})^{2}$$

Linear regression

If we take the derivatives with respect to β₀ and β₁ we will find exactly the same first order conditions as with. For example, with respect to β₁:

 $\sum_{i=1}^{n} x_1 (y_i - \beta_0 - \beta_1 x_1) = 0$

- All the algebraic properties of OLS still hold true here
- The MLE estimate of σ^2 will be biased but we divide by (n-p-1) instead as we saw before
- So what do we gain with MLE?
- We do gain a lot in the understanding of linear regression (more in a little bit)

The regression command again

 The regression command does not use MLE but it does give you the log likelihood

```
sysuse auto, clear
qui reg price weight mpg
ereturn list
scalars
                 e(N) = 74
              e(df m) = 2
              e(df r) = 71
                 e(F) = 14,7398153853841
                e(r2) = .2933891231947529
              e(rmse) = 2514.028573297152
               e(mss) = 186321279.739451
               e(rss) = 448744116.3821706
              e(r2 a) = .27348459145376
               e(11) = -682.8636883111164
              e(11\ 0) = -695.7128688987767
              e(rank) = 3
```

- The log likelihood of the estimated model is stored in e(II). The log likelihood of the null model (with no covariates) is stored in $e(II_0)$.
- From the numbers above $e(II) > e(II_0)$

The regression command again

 Stata uses a formula to go from SSE to log likelihood. Remember, SSE is Stata is stored in the scalar r(rss)

sysuse auto, clear qui reg price weight mpg * Save sample size and SSE local N = e(N) local rss = e(rss) * Use formula local 11 = -0.5*'N'*(ln(2*_pi)*ln('rss'/'N')*1) display %20.6f 'l1' -682.863688 display %20.6f e(l1) -682.863688

• The formula is $II = -0.5N(In(2\pi) + In(\frac{SSE}{N}) + 1)$

Easy MLE in Stata

To estimate in MLE using Stata you need to write a program but Stata now makes it a lot easier (for teaching purposes) with the mlexp command

mlexp (ln(normalden(price, {xb: weight mpg cons}, {sigma}))) initial log likelihood = -<inf> (could not be evaluated) feasible log likelihood = -803.76324 rescale: log likelihood = -729.85758 rescale eq: log likelihood = -697.2346 Iteration 0: log likelihood = -697.2346 Iteration 1. log likelihood = -687.4506 Iteration 2: log likelihood = -682.92425 Iteration 3: log likelihood = -682.86401 Iteration 4: log likelihood = -682.86369 Iteration 5: log likelihood = -682.86369 Maximum likelihood estimation Number of obs = 74 Log likelihood = -682.86369 _____ Coef. Std. Err. z P>|z| [95% Conf. Interval] ----xb weight | 1.746559 .6282189 2.78 0.005 .5152727 2.977846 mpg | -49.51222 84.39157 -0.59 0.557 -214.9167 115.8922 _cons | 1946.069 3523.382 0.55 0.581 -4959.634 8851.771 /sigma 2462.542 202.4197 12.17 0.000 2065.806 2859.277

Almost same

The SEs are slightly different and so is Root MSE. Stata is using the second derivatives to calculate SEs using MLE

. reg price weight mpg

Source	SS	df	MS	Numbe	er of obs	=	74
+				- F(2,	71)	=	14.74
Model	186321280	2	93160639.	9 Prob	> F	=	0.0000
Residual	448744116	71	6320339.6	7 R-squ	lared	=	0.2934
+				- Adjl	R-squared	=	0.2735
Total	635065396	73	8699525.9	7 Root	MSE	=	2514
price	Coef.	Std. Err.	t	P> t	[95% Co	nf.	Interval]
+							
weight	1.746559	.6413538	2.72	0.008	.46773	6	3.025382
mpg	-49.51222	86.15604	-0.57	0.567	-221.302	5	122.278
_cons	1946.069	3597.05	0.54	0.590	-5226.24	5	9118.382

But they are consistent

- The auto dataset has only 74 obs. What about we use the MEPS that has about 15000? (That's really an overkill but just to make the point)
- OLS version:

use heus_mepssa gen lexp = log(reg lexp age fe	exp_tot +1)	.ear					
Source	SS	df	MS	Numbe	r of obs	=	19,386
+-				F(2,	19383)	=	2219.86
Model	35570.262	2	17785.131	Prob	> F	=	0.0000
Residual	155293.552	19,383	8.01184298	R-squ	ared	=	0.1864
+-				Adj R	-squared	=	0.1863
Total	190863.814	19,385	9.8459538	Root	MSE	=	2.8305
lexp		Std. Err.	 t 1			 nf.	Interval]
+-	.0692295	.0011701	59.16 (0.000	. 0669359		.0715231
age female	1.158043	.040891		0.000	1.077893		1.238193
	2.176568	.0603543		0.000	2.058268	-	2.294868
_cons	2.1/0500	.0603543	36.06 (0.000	2.056266	> 	2.294000

Compare to MLE version

mlexp (ln(normalden(lexp, {xb: age female _cons} , {sigma})))

initial:	log likelihood =	- <inf></inf>	(could not be evaluated)
feasible:	log likelihood =	-1539127.9	
rescale:	log likelihood =	-52965.16	
rescale eq:	log likelihood =	-52965.16	
Iteration 0:	log likelihood =	-52965.16	
Iteration 1:	log likelihood =	-48279.366	
Iteration 2:	log likelihood =	-47689.334	
Iteration 3:	log likelihood =	-47676.448	
Iteration 4:	log likelihood =	-47676.408	
Iteration 5:	log likelihood =	-47676.408	

Maximum likelihood estimation

Log likelihood	= -47676.408	В		Number	of obs =	19,386
I	Coef.	Std. Err.	z	P> z	[95% Conf	. Interval]
+-						
xb I						
age	.0692295	.0011701	59.17	0.000	.0669362	.0715227
female	1.158043	.0408878	28.32	0.000	1.077905	1.238182
_cons	2.176568	.0603497	36.07	0.000	2.058285	2.294851
+-						
/sigma	2.830301	.0143739	196.91	0.000	2.802129	2.858473

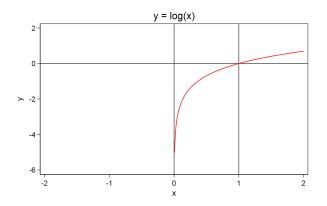
Numerical methods

- Stata uses numerical methods to maximize the likelihood. There are many and some work better than others in some situations. Type "help mle" for the gory details
- A classic one is the Newton-Raphson algorithm
- The idea requires Taylor expansions (a way to approximate nonlinear functions using linear functions)
- The steps are:
 - 1 Make a guess about the parameters, say just one parameter θ_0
 - 2 Approximate the log likelihood function using Taylor series at θ_0 and set it equal to zero (easier to solve because it's a linear function)
 - 3 Find the new θ , say, θ_1 . Check if the log likelihood has improved
 - 4 Repeat until the -2 log likelihood changes by only a small amount, say 0.02
- The idea of using -2 log likelihood < 0.02 is that that amount would not change the statistical significance -2 log likelihood is in the Chi-square scale (more on this in a sec)

Why is the log likelihood function negative?

The likelihood function L(p) is a small number since it's the joint likelihood of observing the outcome values

```
twoway function y = \log(x), range(-2 2) xline(0 1) yline(0) /// color(red) title("y = \log(x)") graph export log.png, replace
```



What do we get using MLE?

- 1) It is clear that we are modeling a **conditional expectation function**: *E*[*Y*|*X*]
- Perhaps this got lost but it's worth repeating. We started with the normal density:

•
$$f(y; \mu, \sigma) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(\frac{-(y_i - \mu)^2}{2\sigma^2}\right)$$

■ We then said that the mean μ is a **function of one or more covariates** x and we made **no assumptions** about the distribution of x:

•
$$f(y; \mu, \sigma) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(\frac{-(y_i - (\beta_0 + \beta_1 x_i))^2}{2\sigma^2}\right)$$

- That's why I said many times that the assumption $\epsilon \sim N(0, \sigma^2)$ is the same as saying that the assumption is $y \sim N(\beta_0 + \beta_1 x, \sigma^2)$, since $\mu = \beta_0 + \beta_1 x$
- Note that with MLE we did not assume anything about the errors. In fact, the errors are not even in the equations. It's a different of deriving the simple linear model

What do we get from MLE?

- 2) It is clear from the start of setting up the problem that we are assuming that Y distributes normal conditional on the values of X. Remember the example of heights for men and women. In some cases, perfectly valid to use a linear model even if the distribution of Y does not look like a normal
- 3) It is clear that we assume that the observations are independent; otherwise, we cannot multiply the densities
- 4) The value of the optimal log likelihood function gives us a measure of the goodness of fit, much like SSR (i.e. the explained part) did. By comparing the log likelihood of alternative models, we will test if the reduced model is adequate like we did with the F test
- 5) The **curvature** of the log likelihood function provides information about the precision of the estimates (i.e. standard errors)

What do we get from MLE?

- 6) MLE is much more general than OLS. You will use MLE for logit, Probit, Poisson, mixture models, survival models. Pretty much all the standard models an applied researcher needs
- 7) Learning to model using likelihood ratio tests is more useful for more type of models than using the SSE for nested models (although you need to use the test command with the robust option in reg)
- 8) AIC and BIC to compare non-nested models are based on the log likelihood function
- Here is a more detailed proof of MLE for the normal: https://www.statlect.com/fundamentals-of-statistics/ normal-distribution-maximum-likelihood

Likelihood ratio test (LRT)

- The null H_0 is that the restricted (constrained) model is adequate
- The alternative H_1 is that the full (unconstrained) model is adequate
- The likelihood ratio test compares the log-likelihoods of both models and can be written as:
- LR = -2[L(RM) L(FM)], where L(RM) is the log-likelihood of the restricted model and L(FM) that of the full model
- Under the null that the restricted model is adequate, the test statistics LR distributes χ^2 with degrees of freedom given by $df = df_{full} df_{restricted}$; that is, the difference in degrees of freedom between the restricted and full models

Likelihood ratio test: sketch of theory

- The theory of LRTs is a bit dense but the intuition is not that difficult to understand
- We could re-write as $LR = -2L(\frac{RM}{FM})$ since $log(\frac{a}{b}) = log(a) log(b)$
- So we are comparing the likelihood of the reduced model to the full model and wondering if the reduced model alone is just fine. Sounds familiar? Not that different from the F test comparing SSEs of nested models
- Keep in mind that the estimated model parameters are those that maximized the value of the likelihood
- The more theoretical part is to figure out how the LRT distributes and under which conditions the LRT is valid (models must be nested)

Recall the F test

- We have $LR = -2L(\frac{RM}{FM})$
- The F test was $F = \frac{[SSE(RM) SSE(FM)]/(p+1-k)}{SSE(FM)/(n-p-1)}$
- Both are using a measure of fit to compare models
- With MLE, we want to know if reaching a higher likelihood is due to chance under the null
- With the F test, we want to know if the additional reduction in the residual variance is due to chance under the null
- The requirement is that models must be nested

Example

Compare the likelihood and other criteria

```
use "GPA1.dta", clear
rename colGPA colgpa
rename hsGPA hsgpa
quietly {
  reg colgpa
  est sto m1
  reg colgpa hsgpa
  est sto m2
  reg colgpa hsgpa skipped
est sto m3
3
est table m1 m2 m3, star stat(r2 r2_a ll bic aic) b(%7.3f)
   Variable | m1 m2
                                 m3
-----
     hsgpa | 0.482*** 0.459***
   skipped |
                              -0.077 **
     _cons | 3.057*** 1.415*** 1.579***
-----+
       r2 | 0.000
                    0.172
                              0.223
     r2_a 0.000 0.166 0.211
      11 | -60.257 -46.963 -42.493
bic | 125.462 103.823 99.832
      aic | 122.513 97.925 90.985
```

legend: * p<0.05; ** p<0.01; *** p<0.001

Example



Irtest m3 m2 Likelihood-ratio test (Assumption: m2 nested in m3)	LR chi2(1) = Prob > chi2 =	8.94 0.0028
. lrtest m3 m1		
Likelihood-ratio test (Assumption: m1 nested in m3)	LR chi2(2) = Prob > chi2 =	35.53 0.0000

 It seems logical that LRT and F-test comparing nested models should be equivalent (asymptotically)

LRT and F-tests

Compare tests

qui reg colgpa est sto mO scalar 110 = e(11)reg colgpa male campus Source | SS df MS Number of obs = 141 -----F(2, 138) = 0.62 Prob > F = 0.5413 Model | .171856209 2 .085928105 138 .139378574 R-squared = 0.0089 Residual | 19.2342432 Adj R-squared = -------0.0055Total | 19.4060994 140 .138614996 Root MSE -.37333 est sto m1 scalar 111 = e(11)1rtest m0 m1 Likelihood-ratio test LR chi2(2) =1 25 (Assumption: m0 nested in m1) Prob > chi2 = 0.5341 * By hand

* By hand di -2*[110 - 111] 1.2542272

■ p-value of both 0.5341 (I chose bad predictors so p-values would be high)

Information criteria to compare models

- A common way to compare models is by using the so-called information criterion. It's a measure that balances bias and variance or accuracy (fit) and simplicity (parsimony)
- The first one proposed is the Akaike Information Criterion (AIC). In the context of the linear model (estimated using OLS), it is:

$$AIC_p = n * ln(\frac{SSE_p}{n}) + 2 * p$$

- p is the number of estimated parameters (including the constant), n is the number of observations, and SSE is the residual sum of squares (recall SST = SSR + SSE)
- The smaller the AIC the better. When is a model going to be better? The larger the sample size, the lower the unexplained variance, the fewer parameters we use
- Usually, we compare models with the same sample size, n, or maybe a small difference in n (a covarite may have some missings)

Log-likelihood and AIC

- That 2 * *p* is a **penalty** much like in adjusted *R*². The more parameters we estimate the better the model has to be to justify the additional estimated parameters
- We just saw the formula linking the log-likehood to SSE: $II = -0.5 * n * [In(2 * \pi) + In(SEE/n] + 1$
- We could solve for SSE and plug in to get AIC as a function of the log-likelihood
- That's one one way to get some intuition on AIC

Information theory

- AIC comes from information theory. Suppose that we know the population process that generated the data but we are using a model to represent that process
- Think of simulating some data and then using imperfect models to represent the data. There will be some information lost
- If we knew the true data generating process, we could find a way to estimate how much information we are losing in each model and thus choose the model that doesn't lose as much information
- But we never know the true model/process. Akaike, in 1974, showed that his information criterion, now called AIC, tells how much more information is lost comparing one model to the other using the estimated log-likelihood

Some caveats

1 Valid asymptotically

- 2 Balance between goodness of fit and parsimony (simplicity). The penalty is incurred to avoid overfitting models because models with more parameters will always fit the data better. If you use as many parameters as data points, model fit is perfect (essentially, the estimated model is a perfect representation of the data)
- **3** Most important: AIC is a relative measure that compares one model to another to choose the one that loses less information. It's not a measure of how good a model is. You may have two bad fitting models (say, super low R^2). AIC helps you choose the best among terrible options

AIC and F-tests

- Note that AIC is similar to the F-test we used to compare nested models. One way of writing the F-test was

 [SSE(RM)-SSE(FM)]/(p+1-k)
 - $F = \frac{[SSE(RM) SSE(FM)]/(p+1-k)}{SSE(FM)/(n-p-1)}$
- Also, the F-test can be written in terms of the R^2 : $F = \frac{(R_p^2 R_q^2)/(p-q)}{(1-R^2p)/(n-p-1)}$
- We are essentially comparing SSE (or *R*²) in relationship to the number of estimated parameters
- AIC is helpful because we can use it to **compare non-nested models**, although there is no statistical test like F test or MLE
- AIC is a relative measure. The absolute value doesn't tell us much; we use it compare related models
- Obviously, it wouldn't make much sense to compare models with different outcomes or different datasets

AIC with log-likelihood

- We also saw that we can use the likelihood ratio test to compare models and that the F-test and the LRT are asymptotically equivalent
- Stata's way to calculate AIC (except in GLM models) is AIC = -2II + 2k
- *k* is the number of parameters, *ll* is the log-likelihood function. 2*k* is the penalty due to the number of parameters; the more parameters, the higher AIC must be to compensate
- Maybe the best way to summarize is that we are essentially comparing the log-likelihood of related models

Bayesian Information Criterion (BIC)

- The more commonly used information criterion is BIC: BIC = -2II + In(n) * k
- Note that we now add ln(n) in the penalty term rather than 2
- If n = 8, ln(8) = 2.079. So in most applications, the BIC penalty is always **higher than the AIC penalty** but it has a limit because the log function "flattens"
- $\ln(50) = 3.91$, $\ln(100) = 4.60$, $\ln(500) = 6.21$, $\ln(1000) = 6.91$, $\ln(2000) = 7.6$
- The more observations we have, the lower the growth of the penalty ln(n) * k

Logit and probit: Latent variable approach

- There is yet another way to derive logit and probit models. A way that connects both into the same framework
- Suppose that there is a latent (unobserved) and continuous variable y* that take values from −∞ to +∞. We assume that the latent variable is a function of covariates X. For simplicity, let's just assume a linear relationship and just one covariate: y_i* = β₀ + β₁x_i + u
- u plays the same role as ϵ in the linear model: a source of random error
- We do not observe the latent variable *y**, we only observe if an event happens or not but whether the event happens depends on the value of the latent variable. We use *y_i* to denote the **observed** variable, which we assume is coded as 1 or 0
- If $y_i * > 0$ then $y_i = 1$. If $y_i * \le 0$ then $y_i = 0$. Note that in this case 0 is a **threshold**
- Think of y* as intelligence and y is whether a person answers a question correctly or not. Or think of the latent variable a measure of disease severity. If it crosses 0 the person dies

Latent variable approach

Because of the way we set up the problem, we can write the probability of y = 1 conditional on the covariate x as:

P(y=1|x) = P(y*>0|x)

- Since we assumed that $y_i^* = \beta_0 + \beta_1 x_i + u$ the above equation becomes $P(y = 1|x) = P(\beta_0 + \beta_1 x_i + u > 0|x) = P(u < [\beta_0 + \beta_1 x_i]|X) = F([\beta_0 + \beta_1 x_i]|x)$
- So the probability of **observing** y = 1 depends on the distribution of u, the error component of the *latent* variable y*. We could calculate that probability if we knew the **cumulative distribution function** F()
- It's obvious that we need to make a strong assumption about the distribution of y* in order to calculate P(y = 1|x) (note that P(y = 0|x) = 1 − P(y = 1|x))

Latent variable approach

- In econometrics this type of model is called index function models (in this case a single index)
- Two common assumptions about *u* are used: *u* distributes either **standard logistic** or **standard normal**
- Both distributions have a mean of 0 and constant variance. In the standard logistic the variance $var(u) = \frac{\pi^2}{3}$. In the standard normal var(u) = 1
- The idea of fixing the variance is not that trivial in the sense that if we don't fix it, then we can't estimate it because we only observe a 0 or 1 and the probability of 1 depends on the sign on *y** but not the scale (variance)
- In other words, we don't have information to estimate var(u) yet we lose nothing by fixing it because P(y = 1|x)does not depend on var(u)

Latent variable approach

- Remember that the cumulative distribution function (cdf) gives you P(X < a). Remember too that to calculate the probability we need to integrate the density f(t) from -∞ to a: ∫^a_{-∞} f(t)dt
- If we assume **standard normal cdf**, our model becomes $P(y = 1|x) = \int_{-\infty}^{\beta_0 + \beta_1 x} \frac{1}{2\pi} e^{(-\frac{t^2}{2})} dt$
- And that's the **probit** model. Note that because we use the cdf, the probability will obviously be constrained between 0 and 1 because since it's a cdf
- If we assume that u distributes **standard logistic** then our model becomes $P(y = 1|x) = \frac{e^{\beta_0 + \beta_1 x}}{1 + e^{\beta_0 + \beta_1 x}}$, which we saw before as the logit transformation
- Remember that there are two different concepts: logistic response function and logistic distribution. The standard logistic cdf happens to have the above formula (the pdf is different)

Estimation

- Estimation is straightforward with MLE. We did it for the logistic model already
- For probit, the likelihood is just like writing P(y = 1|x) above because that's the probability of seeing the data. We need to multiply *n* times and also consider that the probability of 0 is 1 P(y = 1|x). If we take the log, it's a sum
- This is often a source of confusion but remember that the likelihood function is the probability of seeing the data given assumptions about the distribution of the data
- So what is the probability of observing a data point y = 1? It's $P(y = 1|x) = \int_{-\infty}^{\beta_0 + \beta_1 x} \frac{1}{2\pi} e^{(-\frac{t^2}{2})} dt$
- What is the probability of observing a data point y = 0? It's $P(y = 0|x) = 1 \int_{-\infty}^{\beta_0 + \beta_1 x} \frac{1}{2\pi} e^{(-\frac{t^2}{2})} dt$
- For the logit, we would do the same, but we would arrive to the same place as before

Estimation

- We can program the probit model easily with Stata
- I have more examples on my site https://clas.ucdenver.edu/marcelo-perraillon/code-and-topics
- Note below that writing the likelihood makes it obvious that the betas are shifts in the standard normal cdf scale

```
program probit_lf
    version 12
    args todo b lnf
    tempvar xb lj
    mleval 'xb' = 'b'
    * latent variable assumed cumm standard normal
    qui gen double 'lj' = normal('xb') if $ML_y1 == 1
    qui replace 'lj' = normal('xb') if $ML_y1 == 0
    qui mlsum 'lnf' = ln('lj')
end
```

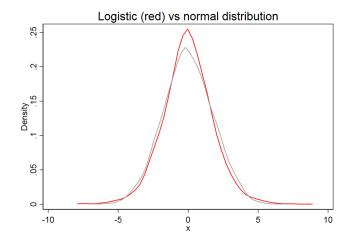
Standard logistic vs standard normal

Does it make much of a difference if we use one vs the other?

```
clear
set seed 123456
set obs 5000
gen u = uniform()
* Simulate logistic distribution
gen 1 = -\ln((1 - u)/u)
sum 1
* Simulated normal with same parameters
gen n = rnormal(r(mean), r(sd))
* Plot.
kdensity 1, bw(0.3) gen(x1 d1)
kdensity n, bw(0.3) gen(xn dn)
line dl xl, sort color(red) || line dn xn, sort ///
   title("Logistic (red) vs normal distribution") ytitle("Density") ///
  xtitle("x") legend(off)
graph export logvsnorm.png, replace
```

Standard logistic vs standard normal normal

Assuming either one as the latent distributions makes little difference



Digression

- Assuming standard normal cdf or logistic are not the only options
- There is the complementary log-log model commonly used in discrete time survival because the exponent of coefficients are hazard rates
- Or the Gumbel model used to model extreme values
- Or the Burr model. Or the Scobit model
- Statistics and econometrics are large fields... Papers must be written, dissertations must be completed
- Sometimes a proposed new method goes to the Journal Article Graveyard. Sometimes they are resurrected 30 years later when somebody discovers that they are perfect for a particular application
- So many ideas and clever people out there. See Greene (2018) for more details

Example

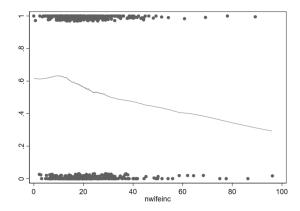
Women's labor force participation (inlf); main predictor is "extra" money in family

bcuse mroz, clear

inlf	=1 if in labor force, 1975							
nwifeinc	(faminc - wage*hours)/1000							
educ	years of schooling							
exper	actual labor mkt exper							
age	woman's age in yrs							
kidslt6	kidslt6 # kids < 6 years							
kidsge6 # kids 6-18								
Variable	1	Obs	Mean	Std. Dev.	Min	Max		
	+-							
i	nlf	753	.5683931	. 4956295	0	1		
nwife	inc	753	20.12896	11.6348	0290575	96		
e	duc	753	12.28685	2.280246	5	17		
ex	per	753	10.63081	8.06913	0	45		
	age	753	42.53785	8.072574	30	60		
kids	1t6	753	.2377158	.523959	0	3		
kids	ge6	753	1.353254	1.319874	0	8		

Labor force participation

The probability of working is decreasing as a function of "extra" income



Writing down the model

We want to estimate the following model:

 $P(inlf_i = 1 | nwifeinc_i) = \Lambda(\beta_0 + \beta_1 nwifeinc_i)$

- By convention (in economics and health economics), when we write capital lambda, Λ(), we imply a logistic model (Λ is not a non-linear function).
 When we write phi, φ(), we imply a probit model
- Write the logistic model this way (no error term!):

$$log(\frac{inlf_i}{1-inlf_i}) = \beta_0 + \beta_1 nwifeinc_i$$

∎ Or

 $logit(inlf_i) = \beta_0 + \beta_1 nwifeinc_i$

• Again, write it like this: $log(\frac{inlf_i}{1-inlf_i}) = \beta_0 + \beta_1 nwifeinc_i$ because this will **match Stata's (or any other statistical package) output**. Remember, we are not directly estimating $P(inlf_i = 1|nwifeinc_i)$

Estimating the model

• So, we will estimate
$$log(\frac{inlf_i}{1-inlf_i}) = \beta_0 + \beta_1 nwifeinc_i$$

logit inlf nwifeinc, nolog

Logistic regres	sion			Number	of obs	=	753
				LR chi2	2(1)	=	10.44
				Prob >	chi2	=	0.0012
Log likelihood = -509.65435				Pseudo	R2	=	0.0101
inlf	Coef.	Std. Err.	z	P> z	[95%	Conf.	Interval]
+-							
nwifeinc	0207569	.0065907	-3.15	0.002	033	6744	0078394
_cons	.6946059	.1521569	4.57	0.000	. 39	6384	.9928279

■ A one thousand increase in "extra" income decreases the log-odds of participating in the labor force by 0.021. And it's statistically significant (p-value = 0.002). Same Wald test as before: -.0207569/.0065907 = -3.1494227. The difference is that the it's not t-student distributed but normally distributed

Overall significance

- The χ² (chi-square) test of the overall significance should look familiar. It compares the current model to the null model (without covariates); the null hypothesis is that all the cofficients in current model are zero
- It's the likelihood ratio test that we have seen before; the equivalent of ANOVA:

```
* LRT
qui logit inlf nwifeinc, nolog
est sto full
qui logit inlf, nolog
est sto redu
lrtest full redu
Likelihood-ratio test LR chi2(1) = 10.44
(Assumption: redu nested in full) Prob > chi2 = 0.0012
```

What about that Pseudo R^2 ?

- We can't partition the variance into explained and unexplained as before so we don't have a nice R² that goes from 0 to 1
- But one way to come up with a measure of fit is to use the (log) likelihood function to compare the current model to the model without any explanatory variable (the null model)
- The formula is: $1 \frac{II_{cm}}{II_{nul}}$, where II_{cm} is the log-likelihood of the current model and II_{nul} is the log-likelihood of the null model
- In other words, adding variables doesn't improve the likelihood. If adding variables improves the likelihood, then the pseudo *R*² will be greater than zero

Pseudo- R^2

Replicate Pseudo R²

```
qui logit inlf muifeinc, nolog
scalar ll_cm = e(ll)
qui logit inlf, nolog
scalar ll_n = e(ll)
di 1 - (ll_cm/ll_n)
.0101362
di "cm: "ll_cm " " "null: " ll_n " " "(ll_cm/ll_n): " (ll_cm/ll_n)
 cm: -509.65435 null: -514.8732 (ll cm/ll n): .9898638
```

- Psuedo R² is **not** a measure of how good the model is at prediction; just how better it fits compared to null model. I don't think that calling it pseudo R² is a good idea
- Big picture: comparing the log-likelihood of models is a way of comparing goodness of fit. If nested, we have the a test (LRT); if not nested, we have BIC or AIC

Not the only pseudo R^2 ?

- Stata uses one version of pseudo *R*² but there are plenty more. Other software may use different metrics
- Long and Freese (2014) have a laundry list of different pseudo R^2 (it's an excellent book, by the way)
- There is the McFadden one, MLE, Cragg and Uhler (also known as Nagelkerke), Efron's, Tjur's... (page 127)
- \blacksquare In any case, none of them have the same meaning as the R^2 in linear regression
- In particular, they **don't mean that predictions are good**. Recall that in linear regression the *R*² is also the square of the correlation between observed and predicted values
- See, context matters a lot

Let's try a different predictor

```
• We will estimate log(\frac{inlf_i}{1-inlf_i}) = \beta_0 + \beta_1 hsp_i, where hsp if education > 12
```

```
gen hsp = 0
replace hsp = 1 if educ > 12 & educ ~= .
logit inlf hsp, nolog
Logistic regression
                                        Number of obs
                                                               753
                                        LR chi2(1)
                                                             15 08
                                        Prob > chi2
                                                            0 0001
Log likelihood = -507.33524
                                        Pseudo R2
                                                             0 0146
       inlf |
                Coef
                       Std Err
                                    7
                                        P>171
                                                 [95% Conf. Interval]
------
              .6504074
                       .1704773
                                  3.82
                                        0.000
                                                 .3162781
                                                           9845368
       hsp |
              0998982
                         086094
                                  1 16
                                        0 246
                                                 - 068843
                                                           2686393
      cons |
```

The log-odds of entering the labor force is 0.65 higher for those with more than high school education compared to those with high-school completed or less than high-school

Odds ratios

- Let's do our usual math to make sense of coefficients. We just estimated the model log(^{inlf_i}/_{1-inlf_i}) = β₀ + β₁hsp_i
- For those with hsp = 1, the model is $log(\frac{inlf_{hsp}}{1-inlf_{hsp}}) = \beta_0 + \beta_1$
- For those with hsp = 0, the model is $log(\frac{inlf_{nohsp}}{1 inlf_{nohsp}}) = \beta_0$
- The difference of the two is $log(\frac{inlf_{hsp}}{1-inlf_{hsp}}) log(\frac{inlf_{nohsp}}{1-inlf_{nohsp}}) = \beta_1$
- Applying the rules of logs: $log(\frac{\frac{inf_{hsp}}{1-inf_{hsp}}}{\frac{inf_{hsp}}{1-inf_{rohsp}}}) = \beta_1$

• Taking
$$e^{()}$$
: $\frac{\frac{infl_{hsp}}{1-infl_{hsp}}}{\frac{infl_{hohsp}}{1-infl_{hohsp}}} = e^{\beta_1}$

Odds ratios

$$rac{rac{inlf_{hsp}}{1-inlf_{hsp}}}{rac{inlf_{nohsp}}{1-inlf_{nohsp}}} = e^{eta_1}$$

And that's the (in)famous odds-ratio

- In our example, $e^{.6504074} = 1.92$. So the odds of entering the labor force is almost twice as high for those with more than high school education compare to those without
- That's the way careful reporters would report this finding. And it's correct. The problem is that we would then interpret this as saying that the probability of entering the labor force is twice as high for those with more than high school
- That interpretation is wrong. A ratio of odds is more often than not far away from the ratio of probabilities

Predictions in the probability scale - logit and probit

- \blacksquare We can easily calculate predictions in the $\ensuremath{\text{\ probability\ scale}}$
- Remember, we are modeling $log(\frac{p}{1-p}) = \beta_0 + \beta_1 X_1 + \dots + \beta_j X_j$
- But we can solve for p:

 $p(\mathbf{X},\beta) = \frac{e^{\beta_0 + \beta_1 X_1 + \dots + \beta_j X_j}}{1 + e^{\beta_0 + \beta_1 X_1 + \dots + \beta_j X_j}} = \frac{1}{1 + e^{-\beta_0 + \beta_1 X_1 + \dots + \beta_j X_j}}$

- That's what Stata does by default when using the postestimation command predict
- Predictions after probit can be a bit confusing, but remember that the coefficients are shifts in the cumulative standard normal (they are z scores), so to calculate p(x, β) we need to calculate a probability given the z score

Calculating predictions after logit and probit models

Make sure you understanding this. Think of predictions after probit as z scores that need to be converted into a probability

logit inlf hsp, nolog

```
* Predictions for logit manually
gen phat_manu = 1/(1*exp(-(_b[_cons] +_b[hsp]*hsp)))
*Same as using the inverse logit function
gen phat_invl = invlogit(_b[_cons] +_b[hsp]*hsp)
* Same as default of predict command
predict phat_pred
```

sum phat_manu phat_invl phat_pred

Variable	Obs	Mean	Std. Dev.	Min	Max		
phat_manu phat_invl phat_pred	753 753 753	.5683931 .5683931	.0694388 .0694388	.5249538 .5249538 .5249538	.6792453		
*probit qui probit inlf hsp, nolog							
<pre>* use inverse normal gem phati_norm = normal(_b[_cons] + _b[hsp]*hsp) predict phati_predprob predict zscore, xb sum phati+ zscore</pre>							
Variable	Obs		Std. Dev.		Max		
phat1_norm phat1_pred [*] b zscore	753	.5683931 .5683931	.0694388 .0694388	.5249538 .5249538 .0625907	.6792453 .6792453		

di normal(0.1760512)

* .5683