

Econometrics Lunch—Simulation-based Estimation

Bernard Salanié

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General references:

- **C. Gouriéroux and A. Monfort (1997)**, *Simulation-Based Econometric Methods*, Oxford University Press.
- **H. van Dijk, A. Monfort and B. Brown (1995)**, *Econometric Inference Using Simulation Techniques*, John Wiley (collected papers).
- **R. Mariano, T. Schuerman and M. Weeks (2000)**, *Simulation-based Inference in Econometrics*, Cambridge University Press (some review papers).

Three generations in econometrics

- 60s: closed-form estimators, mostly OLS on linear models.
- 70s and 80s: the advent of numerical optimization, FIML-GMM-PML.
- 90s and beyond: models too complicated to yield a closed-form criterion \longrightarrow simulate.

The multinomial probit

First example that motivated generation 3.

discrete choice model.

individuals $i = 1, \dots, n$, alternatives $j = 1 \dots, m$

utility of j for i : $U_{ij} = z_{ij}b_j + \varepsilon_{ij}$ where ε_{ij} is $N(0, 1)$, possibly correlations across alternatives j .

$$\begin{aligned}\Pr(i \text{ chooses } j) &= \Pr(\forall k, U_{ij} \geq U_{ik}) \\ &= \Pr(\forall k, \varepsilon_{ij} - \varepsilon_{ik} \geq z_{ik}b_k - z_{ij}b_j)\end{aligned}$$

is an $(m - 1)$ -dimensional integral of the normal distribution, untractable for $m > 5$, which is not unusual for some applications (e.g. transportation).

Second example: model $y_i = g(x_i, \theta, \varepsilon_i, u_i)$

where u_i , the unobserved heterogeneity, has some density h (e.g. panel data with random effects)

Often we can compute in closed form the likelihood conditional on u_i , $f(y_i, x_i, u_i, \theta)$, but not the unconditional likelihood

$$\int f(y_i, x_i, u_i, \theta) h(u_i) du_i$$

Third example: say Tobit with serially correlated latent variable

$$y_{it}^* = x_{it}\beta_0 + \alpha_0 y_{i,t-1}^* + u_{it}$$

Since $y_{i,t-1}^*$ is unobserved, we can't condition on it for estimation
→ the likelihood is a T -dimensional integral.

In these three examples, only simulation will allow us to estimate the model.

The method of simulated moments

Reference: **D. McFadden (1989)**, “A Method of Simulated Moments for Estimation of Discrete Response Models without Numerical Integration”, *Econometrica*, 57, 995-1026.

Here take cross-section data for simplicity.

Often we have a moment condition like

$$E(K(y, x)|x) = k(x, \theta_0)$$

Then normally in GMM we do

$$\min_{\theta} \left\| \frac{1}{n} \sum_{i=1}^n (K(y_i, x_i) - k(x_i, \theta)) \right\|_{\Omega}^2$$

where $\|X\|_{\Omega}^2 = X'\Omega X$, Ω well chosen.

The method of simulated moments, 2

Often k is very hard to compute, as y is driven by a complicated model

$$y = g(x, \theta_0, \varepsilon).$$

(Take the multinomial probit with $K(y_i, x_i) \simeq (i \text{ chooses } j)$, so that $k \simeq P_{ij}$)

then we draw ε^s in the distribution of ε and compute

$$y^s(\theta) = g(x, \theta, \varepsilon^s)$$

and get the MSM estimator by approximating

$$k(x_i, \theta) \text{ with } \frac{1}{S} \sum_{s=1}^S K(y_i^s(\theta), x_i).$$

Is it a good idea?

We just want to approximate an integral, and there are other techniques (see end of slides)

e.g. Gaussian quadrature for some kernel g :

$$\int_0^1 f(x)g(x)dx \equiv \sum_{i=1}^m \omega_i f(x_i)$$

where the (x_i, ω_i) only depend on g .

- much more accurate than Monte Carlo integration for small dimensional integrals
- but in dimensions larger than 3 they become pretty useless.

... and there are ways to improve the precision of Monte Carlo integration in several important cases.

it can be shown that when S is fixed and $T \rightarrow \infty$,

- the MSM estimator is consistent and asymptotically normal
- its asymptotic variance is of the form $(a + \frac{b}{S})$
- so the efficiency loss due to the simulations can be made small by increasing S .

References:

G. Laroque and B. Salanié (1989), “Estimation of Multimarket Fix-Price Models: An Application of Pseudo-Maximum Likelihood Methods”, *Econometrica* , 57, 831-860.

G. Laroque and B. Salanié (1993), “Simulation-Based Estimation of Models with Lagged Latent Variables”, *Journal of Applied Econometrics* , 8, S119-S133.

Take an NLLS example: $E(y|x) = m(x, \theta_0)$.

Often $m(x, \theta)$ is very difficult to compute. So instead we simulate

$$y^s(\theta) = g(x, \varepsilon^s, \theta)$$

Asymptotic properties: although $m^S(x, \theta)$ is unbiased, it enters nonlinearly in the criterion we minimize.

This implies that with fixed S and $n \rightarrow \infty$, the SPML estimators have a bias of order $1/S$.

We can correct the objective function so as to get a consistent estimator for fixed S . To do this, compute the Monte-Carlo variance

$$v^S(x, \theta) = \frac{1}{S-1} \sum_{s=1}^S \left(y^s(\theta) - m^S(x, \theta) \right)^2$$

and minimize over θ

$$\sum_{i=1}^n \left((y_i - m^S(x_i, \theta))^2 - \frac{1}{S} v^S(x_i, \theta) \right)$$

Example: model with unobserved heterogeneity. Then, often

$$f(y, x, \theta, m, \sigma) = \int f(y, x, m + \sigma u, \theta) h(u) du$$

which cannot be integrated (except in some very special cases).
The solution is to draw u^s from h , to let

$$f^S(y, x, \theta, m, \gamma) = \frac{1}{S} \sum_{s=1}^S f(y, x, u^s, \theta, m, \gamma)$$

and to solve

$$\max_{\theta, m, \gamma} \sum_{i=1}^n \log f^S(y, x, \theta, m, \gamma)$$

Other example: models with latent variables (logit, probit, tobit. . .).

often very hard with lagged $y_{i,t-1}^*$ \rightarrow use

$$\frac{\partial \log f}{\partial \theta}(y, x, \theta) = E \left(\frac{\partial \log f^*}{\partial \theta}(y^*, x, \theta) \mid y = h(y^*) \right)$$

where f is the (hard to evaluate) likelihood of the model and f^* is the likelihood of the (unusable) latent-variable model.

so that if we can simulate y^{s*} in $h^{-1}(y)$, we can obtain a SML estimator. This is called the method of simulated scores (MSS).

Asymptotics: f^S is an unbiased simulator of f , but log is nonlinear \rightarrow again, we have a (hopefully small) bias in $1/S$.

Fermanian-Salanié, *Econometric Theory* 2004:

Given a (parametric!) model $y = g(x, \theta, \varepsilon)$,

draw $y^s(x, \theta)$, $s = 1, \dots, S$

and nonparametrically estimate the density:

$$f^S(y, x, \theta) = \frac{1}{Sh} \sum_{s=1}^S K \left(\frac{y - y^s(x, \theta)}{h} \right).$$

Consistent and asymptotically efficient if S goes to infinity and h goes to zero (at the right rates...) as sample size goes to infinity.

Indirect Inference

Reference: **C. Gouriéroux, A. Monfort and E. Renault(1993)**,
“Indirect Inference”, *Journal of Applied Econometrics* , 8,
S85-S118.

Old idea: minimum-distance estimation, resurfaced as asymptotic LS, then (with simulations) as indirect inference.

Assume our model (M) has a complicated likelihood function $f(y, x, \theta)$ that we cannot maximize.

On the other hand, model (M') is reasonably close to model (M) and has a simple likelihood function $f'(y, x, \beta)$ \rightarrow it is easy to estimate.

We could do

$$\max_{\beta} \sum_{i=1}^n \log f'(y_i, x_i, \beta)$$

to estimate (M'). But of course, we get an estimator $\hat{\beta}$ that gives us little information on θ_0 . *Or does it?*

Indirect inference proceeds in three steps:

- 1 for given θ , simulate model (M) to get $y^s(\theta)$
- 2 estimate model (M') on simulated data:

$$\max_{\beta} \sum_{i=1}^n \sum_{s=1}^S \log f'(y^s(\theta), x, \beta)$$

to get an estimator $\hat{\beta}^S(\theta)$

- 3 “calibrate” to make $\hat{\beta}^S(\theta)$ as close to $\hat{\beta}$ as possible, i.e. with some positive definite matrix Ω , do

$$\min_{\theta} \left\| \hat{\beta}^S(\theta) - \hat{\beta} \right\|_{\Omega}^2$$

which gives the indirect inference estimator $\hat{\theta}^S$.

For this to work, we must have $\dim \beta \geq \dim \theta$. Moreover, let b be the “binding function”:

$$b(\theta) = \arg \max_{\beta} E \log f'(y(\theta), x, \beta)$$

$b(\theta)$ is the *pseudo-value*.

Then we need the equation $b(\theta) = b(\theta_0)$ to have the only solution $\theta = \theta_0$:

model (M) is identified from model (M')

The indirect inference estimator $\hat{\theta}^S$ is consistent and asymptotically normal for fixed S , with efficiency loss in $1/S$.

The choice of the auxiliary model matters a great deal. . . just like the choice of instruments in GMM.

Econometrics of finance

The data (say the price of a stock) is generated by a continuous-time process

$$dy_t = g(\theta, y_t)dt + h(\theta, y_t)dW_t$$

where W_t is a Brownian motion. We only observe y_1, \dots, y_T . The problem is how to discretize the model for estimation purposes.

Example, 2

In very rare cases, we can compute the exact discretization. For instance, take the geometric diffusion used by Black-Scholes

$$\frac{dy_t}{y_t} = \mu dt + \sigma dW_t$$

Then Ito's formula gives

$$d(\log y)_t = \left(\mu - \frac{\sigma^2}{2} \right) dt + \sigma dW_t$$

which integrates to

$$\log y_t = \log y_{t-1} + \mu - \frac{\sigma^2}{2} + \sigma \varepsilon_t$$

Usually we can't integrate, so we can use the *naive* discretization

$$y_t = y_{t-1} + g(\beta, y_{t-1}) + h(\beta, y_{t-1})\varepsilon_t$$

as model (M'), which can be estimated through weighted NLLS. This is not the true model, so the 'pseudo-true-value is not θ_0 , but things may work well if it is a (locally) one-to-one function of θ_0 .

Another example: dynamic programming models

Stationary example:

$$V(s_t) = \max_{c_t} (u(c_t, \theta, x_t, \varepsilon_t) + \beta(x_t, \theta)E(V(s_{t+1})|s_t, c_t, \theta)).$$

You observe the three time-series (c_{it}, s_{it}, x_{it}) for several agents, so you “can” estimate nonparametrically (inter alia)

- the transition process $(s_t, c_t, x_t) \longrightarrow s_{t+1}$
- the choice process $(s_t, x_t) \longrightarrow c_t$.

you can then fit a flexible parametric form for these time series processes: $l(s_{t+1}, c_t | s_t, x_t; \beta)$

... and you “can” simulate the dynamic model for any θ

... and use indirect inference to get your estimator for θ

... + overidentification tests since in general $\dim \beta \gg \dim \theta$.

A primer on simulation techniques

the building block usually is a congruential random number generator (RNG).

This starts from an integer “seed” n_0 chosen by the user or the computer and computes recursively

$$n_{m+1} \equiv an_m + b \pmod{c}$$

where a , b and c are well-chosen large numbers.

After a while, this settles into an (almost) uniformly distributed sequence in $[0, c - 1]$. After dividing by c , we get random draws from the uniform $\mathcal{U}[0, 1]$.

A RNG for a distribution F can be obtained by doing $v = F^{-1}(u)$, where u follows $\mathcal{U}[0, 1]$.

Careful: choose a RNG with a large periodicity c .

There are also more clever “direct” tricks to draw from the normal, the Student, the Beta...

E.g. for Student t_n : draw x_i for $i = 1, \dots, n + 1$ in $N(0, 1)$ i.i.d, and generate

$$y = \frac{x_1}{\sqrt{\frac{\sum_{i=2}^{n+1} x_i^2}{n}}}.$$

You want to draw from $f(x)$ and

- you can compute $f(x)$ in every x
- but you cannot draw directly (e.g. F^{-1} is very nasty)
- and you can draw from $g(x)$, another density such that

$$\forall x, \frac{f(x)}{g(x)} \leq \text{some } M \quad (B).$$

Acceptance-rejection algorithm:

- 1 draw x from g
- 2 accept it with proba $f(x)/Mg(x)$
- 3 or go back to step 1.

Acceptance-rejection works in principle, but condition (B) requires g to have thicker tails than f , not so great; and you need to get M about right.

There is a very efficient variant that works by adapting g in several steps:

Adaptive Rejection Sampling

works whenever f is log-concave—which is “very reasonable” unless it has clusters of probability.

Monte Carlo integration

=what we have been doing from the start: approximate $Eh(X) = \int h(x)f(x)dx$ by drawing x_1, \dots, x_n from f and computing

$$\frac{1}{n} \sum_{i=1}^n h(x_i).$$

It has variance Vh/n ; can we do better?

- 1 antithetic variables often work well: use near symmetries in h , e.g:

$$V \left(\frac{h(x) + h(-x)}{2} \right) = \frac{Vh}{2} + \text{cov}(h(x), h(-x)).$$

useful (in this example) when h is “almost odd.”

- 2 importance sampling: draw from g and use

$$\frac{1}{n} \sum_{i=1}^n h(x_i) \frac{f(x_i)}{g(x_i)},$$

clever in theory but often unreliable in practice.

Weighted Monte Carlo Integration

Idea: Riemann sum integration,

$$\int_a^b w(x)dx \simeq \frac{b-a}{n} \sum_{i=1}^n w(a + i(b-a)/n).$$

Bring together with Monte Carlo integration:

- 1 draw the points x_1, \dots, x_n randomly in f
- 2 order them: $x_{(1)} < \dots < x_{(n)}$
- 3 compute

$$\sum_{i=1}^{n-1} h(x_{(i)})f(x_{(i)})(x_{(i+1)} - x_{(i)}).$$

(originally) Bayesian motivation: posterior \leftarrow data + prior

$$f(\theta|y) \simeq f(y|\theta)f(\theta)$$

We are usually concerned with its moments

$$\int \theta^k f(\theta|y) d\theta;$$

but often $\dim \theta$ is large, so we want to do Monte Carlo integration
... provided we can draw from $f(\theta|y)$.

Drawing from the posterior

Often we can evaluate $f(\theta|y)$ at any point, but not draw from it directly.

But there are two MCMC (*Markov Chain Monte Carlo*) tricks:

- Metropolis-Hastings;
- Gibbs sampling.

Say you want to draw from some distribution with pdf $f(x)$;
Choose some density $q(x, y)$ (here, symmetric, for simplicity);
given a draw x^s , draw y from $q(y|x^s)$;
accept it with probability

$$\min \left(\frac{f(y)}{f(x)}, 1 \right)$$

otherwise redraw from $q(y|x^s)$.

After a burn-in phase, draws are as if from f .

If $x = (y, z)$ and you can draw from $f(y|z)$ and $f(z|y)$ but not $f(y, z)$:

given $x^s = (y^s, z^s)$, draw

y^{s+1} from $f(y|z^s)$, z^{s+1} from $f(z|y^s) \longrightarrow x^{s+1}$

again, draws converge to as if from f .

Why is Gibbs sampling so useful?

E.g. in all models with latent Markov chains (Hamilton in macro, volatility models in finance...)

$$y_t = G(s_t, u_t, \theta)$$

We need to draw from $I(s_1, \dots, s_T | y_1, \dots, y_T)$, often very hard, but it is much easier to draw from distributions like

$$I(s_t | s_{-t}, y_1, \dots, y_T).$$

(here if u_t is i.i.d and independent of (s_t) , you just need to condition on (s_{t-1}, y_t, s_{t+1})).