# Lecture 5: Conditional Choice Probabilities (CCPs) Based Methods

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# 1 Introduction

- Finding a fixed point can be computationally demanding.
- Can we somehow skip computing a fixed point when estimating structural parameters?
- Two-step methods do this job.
- A class of two step methods have been proposed in the literature. They are powerful especially in estimating dynamic games.
- To fix the idea, let us discuss the static entry model of Seim (2006).

# 1.1 Example

#### 1.1.1 Model

- Consider a large number of potential entrants:  $N_{\sf max}$
- Parameterization:

$$\Pi(N,S) = S \cdot (\alpha - \gamma(N-1)) - F$$

- $(\alpha, \gamma, F)$  are parameters and (N, S) are observed in each market
- $\epsilon_i$  is private information: Observed by firm i but not other firms or econometrician. Assume iid across players.
- Hence, firm i must form expectation about other firms entering

- Firm i enters if

$$E(\Pi(N,S)|a_i = enter) + \epsilon_i \ge 0$$

or (linearity allows to move through expectations operator)

$$S \cdot (\alpha - \gamma E(N - 1|a_i = enter)) - F + \epsilon_i \ge 0$$

- Assume symmetry

- Let

$$Pr(Entry) = p = G_{\epsilon}(E[\Pi(N, S)|a_i = enter])$$

- Hence, firm i enters if

$$S \cdot (\alpha - \gamma [p \cdot (N_{max} - 1)]) - F + \epsilon_i \ge 0$$

- A Bayesian-Nash Equilibrium is given by a fixed point:

$$p = G_{\epsilon}(S \cdot (\alpha - \gamma[p \cdot (N_{max} - 1)]) - F)$$

- If  $\epsilon_{it}$  iid extreme value:

$$p = \frac{\exp(S \cdot (\alpha - \gamma[p \cdot (N_{max} - 1)]) - F)}{1 + \exp(S \cdot (\alpha - \gamma[p \cdot (N_{max} - 1)]) - F)}$$

- Likelihood contribution of market with N active firms and market demand S:

$$\ell(N, S, p) = \frac{N_{max}!}{(N_{max} - N)!N!} p^{N} (1 - p)^{(N_{max} - N)}$$

#### 1.1.2 Estimation

- Nested fixed point algorithm:
  - 1. Fix parameter vector  $\theta = (\alpha, \gamma, F)$
  - 2. Find fixed point in

$$p_m = G_{\epsilon}(S_m \cdot (\alpha - \gamma [p_m \cdot (N_{max} - 1)]) - F)$$

for every market m = 1...., M, yielding  $p_m(\theta)$ 

3. Evaluate log-likelihood

$$L_M(\theta) = \frac{1}{M} \sum_{m=1}^{M} \log(\ell(N_m, S_m, p_m(\theta)))$$

4. Use numerical algorithm to find  $\theta$  that maximizes  $L_M(\theta)$ 

- Alternatively, we could form an moment estimator:
  - 1. Estimate probability of entry in each market

$$\hat{p}_m = \frac{N_m}{N_{\mathsf{max}}}$$

2. Moment condition:

$$E[\hat{p}_m - G_{\epsilon} (S_m \cdot (\alpha - \gamma [\hat{p}_m \cdot (N_{\mathsf{max}} - 1)]) - F) | Z_m] = 0$$

where  $Z_m$  is a vector of exogenous variables (e.g.,  $S_m$ ). The sample analogues is

$$rac{1}{M}\sum_{m=1}^{M} Z_m \otimes \left[\hat{p}_m - G_\epsilon(S_m \cdot (lpha - \delta[\hat{p}_m \cdot (N_{\sf max} - 1)]) - F)
ight] = 0$$

- Loosely speaking, two-step methods let the data tell us what agents did in the first stage, and choose structural parameters to explain why they did so in the second stage.

- Advantage of two-step estimators:
  - No need to solve for equilibrium for each trial of parameter vector
  - Avoid problem of multiple equilibria
    - Even with multiplicity, we can calculate the equilibrium probability in each market when  $N_{\text{max}}$  is large ("picked by the data") as in step 1 above (see social interaction literature, e.g., Brock and Durlauf, 2001)
    - When  $N_{\text{max}}$  is small, we need to pool different markets
    - Either focusing on unique outcomes or specifying an equilibrium selection rule is needed

# 2 CCP Methods in Dynamic Models

- Can we apply the same idea to our Rust-type model?
- Recall our Bellman equation

$$V(x_t, \varepsilon_t, \theta) = \max_{i \in D(x_t)} \{ u(x_t, i, \theta) + \varepsilon_t(i) + \beta E[V(x_{t+1}, \varepsilon_{t+1}, \theta) | x_t, \varepsilon_t, i] \}$$

- Recall: under CI, the integrated value function formulation in our Rust-type model is

$$V(x_t, \theta) = \mathcal{P}_{x_t} \left[ u(x_t, 1, \theta) + \tilde{\varepsilon}_1^{\mathcal{P}_{x_t}} + \beta E[V(x_{t+1}, \theta) | x_t, 1] \right]$$
$$+ (1 - \mathcal{P}_{x_t}) \left[ u(x_t, 0, \theta) + \tilde{\varepsilon}_0^{\mathcal{P}_{x_t}} + \beta E[V(x_{t+1}, \theta) | x_t, 0] \right]$$

where  $\widetilde{\varepsilon}_{i}^{\mathcal{P}_{x_{t}}} = E[\varepsilon_{t}(i)|i]$  and  $\mathcal{P}_{x_{t}} \equiv \Pr(i_{t} = 1|x_{t}, \theta)$ .

- In matrix notation,

$$\mathbf{V} = \mathbf{\Pi}^{\mathcal{P}} + \beta G^{\mathcal{P},p} \mathbf{V}$$

where the j-th element of  $\Pi$  is

$$\Pi_j^{\mathcal{P}} = \mathcal{P}_{x_t}(u(x_t, 1, \theta) + \widetilde{\varepsilon}_1^{\mathcal{P}_{x_t}}) + (1 - \mathcal{P}_{x_t})(u(x_t, 0, \theta) + \widetilde{\varepsilon}_0^{\mathcal{P}_{x_t}})$$

with  $x_t = x_t(j)$ , and (i, j) element of  $G^{\mathcal{P}, p}$  is

$$G_{ij}^{\mathcal{P},p} = \mathcal{P}_x p(x'|\mathbf{0},\mathbf{1},\theta) + (\mathbf{1} - \mathcal{P}_x)p(x'|x,\mathbf{0},\theta)$$

with  $x = x_t(i)$  and  $x' = x_t(j)$ .

- We get

$$\mathbf{V} = (I - \beta G^{\mathcal{P}, p})^{-1} \Pi^{\mathcal{P}}. \tag{1}$$

- Notice that RHS of equation (1) can be computed from  $\mathcal{P}_{x_t} \equiv \Pr(i_t = 1 | x_t, \theta)$  and  $p(x_{t+1} | x_t, i_t, \theta)$ .

- The basic idea of CCP-based methods is to estimate  $\Pr(i_t = 1|x_t, \theta)$  and  $p(x_{t+1}|x_t, i_t, \theta)$  directly from the data in the first stage, and then estimate structural parameters in the second stage.
- The first two-step method is proposed by Hotz and Miller (1993)
- Notice a big difference from the nested fixed point algorithm: we skip solving a fixed point problem.
- From now, let's assume we observe N individual over T time periods. We use n and t as a subscript to index individual and period, respectively: data  $\{i_{nt}, x_{nt}: n=1,...,N; t=1,...,T\}$
- Typically, T is relatively short; asymptotics is in terms of N.

- Suppose x has a finite support. Then we can use

$$\widehat{\mathcal{P}}_{x} = \frac{\sum_{n=1}^{N} \sum_{t=1}^{T} \mathbf{1}(i_{nt} = 1, x_{nt} = x)}{\sum_{n=1}^{N} \sum_{t=1}^{T} \mathbf{1}(x_{nt} = x)}$$

and

$$\widehat{p}(x'|x,i,\theta) = \frac{\sum_{n=1}^{N} \sum_{t=1}^{T-1} \mathbf{1}(x_{nt+1} = x', x_{nt} = x, i_{nt} = i)}{\sum_{n=1}^{N} \sum_{t=1}^{T-1} \mathbf{1}(x_{nt} = x, i_{nt} = i)}$$

Using these, we can compute

$$\widehat{\mathbf{V}} = (I - \beta G^{\widehat{\mathcal{P}}, \widehat{p}})^{-1} \Pi^{\widehat{\mathcal{P}}}.$$

- Then,

$$\Pr(i_{nt} = 1 | x_{nt}, \theta) = \Pr\left(\widehat{\bar{V}}^{1}(x_{nt}, \theta) + \varepsilon_{1nt} \ge \widehat{\bar{V}}^{0}(x_{nt}, \theta) + \varepsilon_{0nt}\right) \quad (2)$$

where

$$\widehat{\overline{V}^0}(x_{nt},\theta) = u(x_{nt},0,\theta) + \beta E[\widehat{V}(x_{nt+1},\theta)|x_{nt},i_{nt}=0,\theta]$$

$$\widehat{\overline{V}^1}(x_{nt},\theta) = u(x_{nt},1,\theta) + \beta E[\widehat{V}(x_{nt+1},\theta)|x_{nt},i_{nt}=1,\theta].$$

- We can use (2) to form the likelihood for the choices in the data.
- What we do not know are  $u(x_{nt}, 0, \theta)$  and  $u(x_{nt}, 1, \theta)$  only. The parameters in these functions are estimated by MLE.
- Identification of these structural parameters is discussed later. In general, the discount factor and the distribution of  $\varepsilon$  should be assumed known.
- Let's make things more formal and general. Now we have K alternatives, and let the choice be denoted by  $d\in\{1,...,K\}$  .
- For simplicity, suppose  $p(x_{t+1}|x_t, d_t, \theta)$  is known, so we focus on choice probabilities.

- Define

$$egin{aligned} \mathsf{Pr}(d_{nt} = d | x_{nt}, heta) &\equiv \Psi\left(d | x_{nt}; \mathcal{P}, heta
ight) \ &= \mathsf{Pr}\left(d = \mathsf{arg}\max_{d' \in \{1, ..., K\}} \left\{\widehat{ar{V}}^{d'}\left(x_{nt}, heta
ight) + arepsilon_{d'nt}
ight\}
ight) \end{aligned}$$

- Assume (a) the true parameters  $\theta_0$  is identified, and (b) the observations are independent across individuals and  $\Pr(x_{nt} = x) > 0$  for all x in  $\mathbf{X}$ 

- Pseudo likelihood function

$$Q_N\left(\theta,\mathcal{P}
ight) = rac{1}{N} \sum_{n=1}^{N} \sum_{t=1}^{T} \ln \Psi\left(d_{nt}|x_{nt};\mathcal{P},\theta
ight).$$

- The MLE for the nested-fixed point algorithm is

$$\hat{ heta}_{MLE} = rg\max_{ heta \in \Theta} Q_N\left( heta, \mathcal{P}
ight) ext{ s.t. } \mathcal{P} = \Psi\left( heta, \mathcal{P}
ight).$$

- Finding this is costly because for any  $\theta$ , one has to find  $\mathcal{P}$  as a solution to the fixed point problem.

#### 2.1 Pseudo Maximum Likelihood Estimation

- Suppose that we know the population probabilities  $\mathcal{P}^0$  and consider

$$\hat{\theta} \equiv \arg\max_{\theta \in \Theta} Q_N\left(\theta, \mathcal{P}^0\right).$$

- Under standard regularity conditions, this estimator is root-N consistent and asymptotically normal.
- A feasible alternative is

$$\hat{\theta}_{2S} \equiv \arg\max_{\theta \in \Theta} Q_N(\theta, \hat{\mathcal{P}}^0)$$

where  $\hat{\mathcal{P}}^0$  is agent's choice probabilities estimated non-parametrically.

**Result** Suppose (i) Assumptions (a)-(b) hold, (ii)  $\Psi(\theta, \mathcal{P})$  is twice continuously differentiable, (iii)  $\Theta$  is a compact set, (iv)  $\theta^0 \in int(\Theta)$ , and (v) let  $\hat{\mathcal{P}}^0$  be an estimator of  $\mathcal{P}^0$  such that  $\sqrt{N}(\hat{\mathcal{P}}^0 - \mathcal{P}^0) \to_d N(0, \Sigma)$ . Then,  $\sqrt{N}(\hat{\theta}_{2S} - \theta^0) \to_d N(0, V_{2S})$ , where

$$V_{2S} = \Omega_{\theta\theta}^{-1} + \Omega_{\theta\theta}^{-1} \Omega_{\theta\mathcal{P}} \Sigma \Omega_{\theta\mathcal{P}}' \Omega_{\theta\theta}^{-1}$$
 and  $\Omega_{\theta\theta} \equiv E\left(\{\nabla_{\theta}s_n\} \{\nabla_{\theta}s_n\}'\right)$ , with  $s_n \equiv \sum_t \ln \Psi\left(d_{nt}|x_{nt}; \mathcal{P}^0, \theta^0\right)$  and  $\Omega_{\theta\mathcal{P}} \equiv E\left(\{\nabla_{\theta}s_n\} \{\nabla_{\mathcal{P}}s_n\}'\right)$ .

- Drawbacks of the two-step PML: the nonparametric estimator of  $\mathcal{P}^0$  can be very imprecise; and it cannot deal with unobservable individual type.

### 2.2 Nested Pseudo Likelihood Method

- Remember our problem is

$$\hat{\theta} = \arg\max_{\theta \in \Theta} Q\left(\theta, \hat{\mathcal{P}}\right).$$

- A recursive extension of the two-step PML estimator: Aguirregabiria and Mira (2002, 2007)
- Take  $\hat{\mathcal{P}}_0$ , an initial guess of the vector of (possibly non-consistent) agent's choice probabilities.
- A sequence of estimators  $\{\hat{\theta}_K: K \geq 1\}$  defined as

$$\hat{\theta}_K = \arg\max_{\theta \in \Theta} Q(\theta, \hat{\mathcal{P}}_{K-1}) \tag{3}$$

and the probabilities  $\{\hat{\mathcal{P}}_K: K \geq 1\}$  are obtained recursively as

$$\hat{\mathcal{P}}_K = \Psi(\hat{\mathcal{P}}_{K-1}, \hat{\theta}_K). \tag{4}$$

- If this sequence converges, its limit  $(\hat{\theta}, \hat{\mathcal{P}})$  satisfies the following two properties:  $\hat{\theta}$  maximizes the pseudo likelihood  $Q_N(\theta, \hat{\mathcal{P}})$  and  $\hat{\mathcal{P}} = \Psi(\hat{\mathcal{P}}, \hat{\theta})$ . Call any pair  $(\theta, \mathcal{P})$  that satisfies these properties a NPL fixed point.
- Let  $\mathcal{Y}_N$  be the set of NPL fixed points. Then the NPL estimator is defined as

$$(\hat{\theta}_{NPL}, \hat{\mathcal{P}}_{NPL}) = \arg\max_{(\theta, \mathcal{P}) \in \mathcal{Y}_N} Q_N(\theta, \mathcal{P}).$$

- Aguirregabiria and Mira (2007) show that NPL is asymptotically more efficient than the infeasible PML. See also Pesendorfer and Schmidt-Dengler (2010) and Kasahara and Shimotsu (2012).

#### 2.3 Forward Simulation

- Forward simulation is one attractive alternative.
- The Hotz-Miller (1993), Hotz-Miller-Sanders-Smith (1994) idea.
- Let us come back to the simple Rust-type model.
- Recall that under the Extreme value distribution assumption, the choice probability is

$$\Pr(i_t = 1|x_t, \theta) = \frac{\exp(\bar{V}^1(x_t, \theta))}{\exp(\bar{V}^0(x_t, \theta)) + \exp(\bar{V}^1(x_t, \theta))}$$

which implies:

$$\frac{\Pr(i_t = 1|x_t, \theta)}{\Pr(i_t = 0|x_t, \theta)} = \frac{\exp(\bar{V}^1(x_t, \theta))}{\exp(\bar{V}^0(x_t, \theta))}$$

or taking logs:

$$\ln(\Pr(i_t = 1|x_t, \theta)) - \ln(\Pr(i_t = 0|x_t, \theta)) = \bar{V}^1(x_t, \theta) - \bar{V}^0(x_t, \theta)$$

- Hence the choice probabilities (which we can estimate from the data) give us an estimate of the difference in alternative-specific value functions.
- This result (the difference in alternative-specific value functions is uniquely pinned down as a function of CCPs) holds for a wider class of error distributions; "Hotz-Miller inversion".
- Let  $\hat{P}(i, x_t)$  be an estimate of choice probabilities and  $\hat{p}(x_{t+1}|x_t, i_t)$  be an estimate of the transition probability.

- Algorithm:
  - 1. Estimate

$$\widehat{\bar{V}^1}(x_t) - \widehat{\bar{V}^0}(x_t) = \ln(\hat{\mathcal{P}}(1, x_t)) - \ln(\hat{\mathcal{P}}(0, x_t))$$

(I took out  $\theta$  to emphasize that this estimate does not depend on structural parameters; directly recovered from the data)

- 2. Draw S paths of T-length of random shocks  $[(\varepsilon_{0,1}^s, \varepsilon_{1,1}^s), (\varepsilon_{0,2}^s, \varepsilon_{1,2}^s), ..., (\varepsilon_{0,T}^s, \varepsilon_{1,T}^s)]$
- 3. Given shock  $(\varepsilon_{0,t}, \varepsilon_{1,t})$  and  $x_t$  we can find optimal choice because  $i_t=1$  if

$$\widehat{\overline{V}^{1}}(x_{t}) - \widehat{\overline{V}^{0}}(x_{t}) > \varepsilon_{0,t}^{s} - \varepsilon_{1,t}^{s}$$

and  $i_t = 0$ , otherwise.

4. Simulate S paths  $(\tilde{x}_t^s, \tilde{\imath}_t^s)_{t=1}^T$  and corresponding choice specific value func-

tions for given parameters  $(\theta_1, R)$ :

$$\widetilde{V}^{1}(x_{t}, \theta_{1}, R) = -R - c(\mathbf{0}, \theta_{1})$$

$$+ \frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \left( u(\tilde{x}_{t}^{s}, \tilde{\imath}_{t}^{s}, \theta_{1}, R) + \varepsilon_{\tilde{\imath}_{t}^{s}, t}^{s} \right)$$

$$\widetilde{V}^{0}(x_{t}, \theta_{1}, R) = -c(x_{t}, \theta_{1})$$

$$+ \frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \left( u(\tilde{x}_{t}^{s}, \tilde{\imath}_{t}^{s}, \theta_{1}, R) + \varepsilon_{\tilde{\imath}_{t}^{s}, t}^{s} \right)$$

and calculate moments:

$$\frac{1}{\#X} \sum_{x_t \in X} Z_t \otimes \left( \begin{array}{c} \left[ \widetilde{\bar{V}^1}(x_t, \theta_1, R) - \widetilde{\bar{V}^0}(x_t, \theta_1, R) \right] \\ - \left[ \widehat{\bar{V}^1}(x_t) - \widehat{\bar{V}^0}(x_t) \right] \end{array} \right)$$

5. Repeat step 4 until the above is as close to zero as possible (remember that steps 1-3 are fixed in the parameter search

- If we need to perform this forward simulation for any heta, the computation burden may be heavy
- One way to significantly reduce the computation burden is to assume

$$u(x, i, \theta) + \varepsilon_{it} = \psi(x, i, \varepsilon_{it}) \cdot \theta$$

where  $\psi\left(x,i,arepsilon_{i}
ight)$  is a vector of basis functions which does NOT depend on heta

- Sometimes, this comes with no cost (no additional assumption). Consider again our Rust-type model:

$$u\left(x,i, heta
ight) = \left\{ egin{array}{ll} - heta_1 x & ext{if } i=0 \ -R & ext{if } i=1 \end{array} 
ight.$$

- Then, for example, the continuation payoff in  $\widetilde{\bar{V}^1}(x_t, heta_1, R)$  can be written as

$$\frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \left( u(\tilde{x}_{t}^{s}, \tilde{\imath}_{t}^{s}, \theta_{1}, R) + \varepsilon_{\tilde{\imath}_{t}^{s}, t}^{s} \right)$$

$$= \frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \left[ (-\theta_{1} \tilde{x}_{t}^{s}) \mathbf{1}(\tilde{\imath}_{t}^{s} = 0) + (-R) \mathbf{1}(\tilde{\imath}_{t}^{s} = 1) + \varepsilon_{\tilde{\imath}_{t}^{s}, t}^{s} \right]$$

$$= -\theta_{1} \frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \tilde{x}_{t}^{s} \mathbf{1}(\tilde{\imath}_{t}^{s} = 0) - R \frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \mathbf{1}(\tilde{\imath}_{t}^{s} = 1) + \frac{1}{S} \sum_{s=1}^{S} \sum_{t=1}^{T} \beta^{t} \varepsilon_{\tilde{\imath}_{t}^{s}, t}^{s}$$

$$= -\theta_{1} A - RB + C$$

- (A,B,C) do not depend on structural parameters, so for any strategy profile, given by  $(\tilde{x}_t^s, \tilde{\imath}_t^s)_{t=1}^T$ , one can use the forward simulation procedure only once to estimate (A,B,C) and then obtain  $\widetilde{\bar{V}^1}(x_t,\theta_1,R)$  and  $\widetilde{\bar{V}^0}(x_t,\theta_1,R)$  easily for any value of  $\theta$ .

#### - Comments:

- V(x) should be computed only for  $x \in \widetilde{\mathbf{X}}$ , where  $\widetilde{\mathbf{X}}$  is the set of states that are observed in the data and states that are reachable within one time period from each of the observed states. Why?
- But you still have to know CCPs for all the states. Why?
- ullet In practice, we cannot compute an infinite sum. How do we choose T?
- Not exact.

# 2.4 Absorbing States

Another alternative is to exploit absorbing states.

- To explain how absorbing states can ease CCP estimation in general, let us use a single agent optimal stopping problem; continue (i = 0) or stop (i = 1).
- The per-period payoff function is denoted by u(x, i) and we use additively separable payoff shocks  $\varepsilon = (\varepsilon_0, \varepsilon_1)$ . The value function is

$$V\left(x,\varepsilon,\theta\right) = \max_{i\in\{0,1\}} \left\{ u\left(x,i,\theta\right) + \varepsilon_i + \beta E\left[V\left(x',\varepsilon',\theta\right)|x,\varepsilon,i\right] \right\}.$$

Under the conditional independence assumption, alternative-specific value functions (net of  $\varepsilon$ ) are

$$\bar{V}^{0}(x,\theta) = u(x,0,\theta) + \beta E \left[ V(x',\varepsilon',\theta)|x,0 \right]$$
$$\bar{V}^{1}(x,\theta) = u(x,1,\theta) + \beta E \left[ V(x',\varepsilon',\theta)|x,1 \right]$$

- Using this, as in Rust (1987)'s model, the value is

$$V(x,\varepsilon,\theta) = \max_{i\in\{0,1\}} \left\{ \bar{V}^0(x,\theta) + \varepsilon_0, \bar{V}^1(x,\theta) + \varepsilon_1 \right\}.$$

and

$$\bar{V}^{0}(x,\theta) = u(x,0,\theta) + \beta E \left[ \max_{i \in \{0,1\}} \left\{ \bar{V}^{0}(x',\theta) + \varepsilon'_{0}, \bar{V}^{1}(x',\theta) + \varepsilon'_{1} \right\} | x, 0 \right]$$

$$\bar{V}^{1}(x,\theta) = u(x,1,\theta) + \beta E \left[ \max_{i \in \{0,1\}} \left\{ \bar{V}^{0}(x',\theta) + \varepsilon'_{0}, \bar{V}^{1}(x',\theta) + \varepsilon'_{1} \right\} | x, 1 \right]$$

- Assuming  $\varepsilon$  follows the type I extreme value distribution,

$$\bar{V}^{0}(x,\theta) = u(x,0,\theta) + \beta \sum_{x' \in \mathbf{X}} \left[ \mu + \ln \left\{ e^{\bar{V}^{0}(x',\theta)} + e^{\bar{V}^{1}(x',\theta)} \right\} \right] g(x',0,x,\theta)$$

$$(5)$$

$$\bar{V}^{1}(x,\theta) = u(x,1,\theta) + \beta \sum_{x' \in \mathbf{X}} \left[ \mu + \ln \left\{ e^{\bar{V}^{0}(x',\theta)} + e^{\bar{V}^{1}(x',\theta)} \right\} \right] g(x',1,x,\theta)$$

(6)

and the choice probabilities are

$$\Pr(i = 0|x, \theta) = \frac{e^{\bar{V}^{0}(x, \theta)}}{e^{\bar{V}^{0}(x, \theta)} + e^{\bar{V}^{1}(x, \theta)}}$$

$$\Pr(i = 1|x, \theta) = \frac{e^{\bar{V}^{0}(x, \theta)} + e^{\bar{V}^{1}(x, \theta)}}{e^{\bar{V}^{0}(x, \theta)} + e^{\bar{V}^{1}(x, \theta)}}.$$
(8)

$$\Pr(i = 1|x, \theta) = \frac{e^{V^{1}(x, \theta)}}{e^{\bar{V}^{0}(x, \theta)} + e^{\bar{V}^{1}(x, \theta)}}.$$
 (8)

- Now assume that i=1 is a terminating action (that leads to an absorbing state, such as exit), so  $u(x, 1, \theta) = u(x, \theta) + \phi$  and  $u(x, 0, \theta) = u(x, \theta)$ . Then,

$$\bar{V}^{1}(x,\theta) = u(x,\theta) + \phi, \tag{9}$$

where  $\phi$  is the value of the terminating action (e.g., exit value)

- Rearrange (8) to get

$$e^{\bar{V}^{0}(x,\theta)} + e^{\bar{V}^{1}(x,\theta)} = \frac{e^{\bar{V}^{1}(x,\theta)}}{\operatorname{Pr}(i=1|x,\theta)}$$

and plug this in into (5) to have

$$\bar{V}^{0}(x,\theta) = u(x,\theta) + \beta \sum_{x' \in \mathbf{X}} \left[ \mu + \bar{V}^{1}(x',\theta) - \ln \Pr\left(i = 1 | x',\theta\right) \right] g(x',0,x,\theta)$$
(10)

- This expression itself does not look very helpful. But once you use (9), the RHS of (10) is written only in terms of functions of parameters  $(u \text{ and } \phi)$  and something that we can recover from the data  $(Pr\ (i=1|x',\theta) \text{ and } g)$ .
- In other words, in the presence of a terminating action, the one-period-ahead probability of observing such an action can be used to cut off the future dependence, and so the value can be computed.
- This is a special case of finite dependence in Altug and Miller (1998) and Arcidiacono and Miller (2011).

#### 2.5 Practical Issues

- Two step methods are a powerful tool: one does not need to compute a fixed point even once to estimate parameters.
- First-stage estimates may be imprecise: the PML estimator of Aguirregabiria and Mira (2002, 2007) could be helpful.
- The first stage CCP should be consistently estimated. When is it inconsistent?
  - when a researcher pools different markets in the presence of multiple equilibria or unobserved heterogeneity: to test, see Otsu, Pesendorfer, and Takahashi (2016).
  - two-step method allowing for unobserved heterogeneity: See Arcidiacono and Miller (2011) and Otsu, Pesendorfer, Sasaki, and Takahashi (2019).

- Specification of the first stage is up to the researcher; "flexibility-feasibility trade-off"
  - In theory, first stage estimates should be non-parametric.
  - In practice, one may need to impose some functional form assumption.
- What to do with typical data (short T)?
- In non-stationary environments (or the industry in question is still in a transition process during the sample period), some state is rarely observed. We may have to rely on extrapolation.
- What if there are several states that are not observed in the data at all?
  - Interpolation?
  - Extrapolation? See Hu, Sasaki and Takahashi (2016).