

Work after

- 1) I wasn't computing the TC correctly b/c in the last H periods, should feed in $hx^j s_t$ for shock $\{s_t\}_{t-H}^T$.
- 2) If I don't impose TC, should impose TR w/a res, or w/o a res but then have a res in A1 or A2.
- 3) Should check the false thing.

2) Adding TR as a residual eq doesn't change much at all for $\{x, i\}_t$! (stopped prematurely.)
Making the norm the loss concept however gets me much closer to the Taylor-rule!

Adding TR as a res did bring things closer to the TR-outcome for $\{x, i\}_t$ as input. Here the norm doesn't make a big diff. (stopped prematurely.)

Adding TR for $\{i_t\}$ as only input. (local min but VERY unstable.) For norm unstable too, stopped premat.
But it no longer blows up in my face :)

1) Computing the and TC w/ correct $E(\cdot)$ doesn't change things. Does the norm? It finds a local min, but a strange one. (Both for $\{i\}$ only.)

For $\{x_i, i\}$, norm or no norm makes no diff, but I'm in general much closer to TR than before, if I initialize there. If not, then further.

For $\{\pi_i, x_i, i\}$ as input, I get the same as before.
 Not if I initialize at TR, I get a strange local min.
 With norm too.
 at norm almost the same.

↳ Either the method is unstable or I'm not quite doing it correctly
→ Tomorrow: do fsolve thing

✓. Fix H

8 April 2020

✓. Fix too long fightles

✓. Fix f solve for TR & RE-TC \rightarrow fixed! (resids are not defined for $t=1$)

- Stting is f' up w/ the anm-TC.

\hookrightarrow or: I don't know if the sol. procedure is unstable, or if I'm doing sthg wrong.

Back to VFI: Eric Sims does it by surpin' plug in the constraint to rewrite the problem as a fct of the future state:

$$\begin{aligned} \text{We had: } V^{l+1}(k_t) &= \max_{c_t} \ln(c_t) + \beta V^l(k_{t+1}) \\ \text{s.t. } k_t^\alpha - c_t &\geq k_{t+1} \\ \rightarrow c_t &= k_t^\alpha - k_{t+1} \end{aligned}$$

$$V^{l+1}(k_t) = \max_{k_{t+1}} \ln(k_t^\alpha - k_{t+1}) + \beta V^l(k_{t+1})$$

Rewrite in terms of k, k'

$$V^{l+1}(k) = \max_{k'} \ln(k^\alpha - k') + \beta V^l(k')$$

Ok that's fine but I'm idiot - we need the grid for k_t nonetheless! In the nonstochastic world, we just need no evaluation of expectations and possibly no marker chains.

Btw, Eric Sims calls the choice of k' the policy π . But I guess that's isomorphic w/ choosing c_t .

I've found Collard's value function iteration notes.
(value-function-collard-lectnotes.pdf)

→ it seems to suggest that you use the grid for k and k'

for $i = 1:n_{\text{grid}}$

$k_+ = k_{\text{grid}}(i)$

for $j = 1:n_{\text{grid}}$ * there's a complication, but ignore for now.

$k_{++} = k_{\text{grid}}(j)$

$$c(i, j) = c(k_+, k_{++}) \rightarrow u(i, j) = u(c(i, j))$$

$n_{\text{grid}} \times 1$

end

$$j^* = \max[u(c(i, j)), 2]$$

$$k_{++}^* = k_{\text{grid}}(j^*)$$

$$v_{1 \times 1}^{\text{new}}(i) = u(c(i, j^*)) + \beta \cdot v_{1 \times 1}(j^*)$$

end

$$\text{crit} = \max(\text{abs}(v^{\text{new}} - v))$$

Actually Collard does this

$c(i, :)$ is c when $k_+ = k_{\text{grid}}(i)$, and for all k_{++}
 $1 \times n_{\text{grid}}$
 $\rightarrow u(i, :) = u(c(i, :))$ $1 \times n_{\text{grid}}$

$$[v^{\text{new}}, j^*] = \max \left[u(c(i, :)) + \beta v(:) \right]$$

$n_{\text{grid}} \times 1$

* the additional complication is that $c_t, k_{t+1} \geq 0$

$$c_t \geq 0 \Rightarrow c_t = k_t^\alpha - k_{t+1} \geq 0$$

$$k_{t+1} \geq 0 \Rightarrow \boxed{k_t^\alpha > k_{t+1} \geq 0}$$

The grid makes sure that $k_{t+1} \geq 0$, but I need to check that $k_{t+1} \text{ Brt} > k_t^\alpha(i)$ for each i .

Colvard's VFI notes are helpful b/c 9 April 2020

they clarify what the maximization means and what $v(k_{t+1})$ means. $v^{\text{old}}(k_{t+1})$ just means

$v^{\text{old}}(\cdot)$ for all values of the k -grid (k_{t+1})

Similarly, you evaluate $c(k_{t+1}(i), \cdot)$ for all values of the k -grid (k_{t+1}).

The maximization then is just to choose the index

$$j^* \text{ that } \max U(c(k_{t+1}(i), \cdot) + \beta v^{\text{old}}(\cdot),$$

and that for each i . So $k_{t+1}(i) = k_{\text{grid}}(j)$.

$$\text{and } v^{\text{new}}(i) = U(c(k_{t+1}(i), j^*) + \beta v^{\text{old}}(j^*)$$

FSOLVE/FMINCON

↳ could try to attack from several, or smart initialization points.

A potential problem for the anchoring TC:
overparameterized?

I input $T+H = 60$ periods, but I can evaluate only T residuals.

① Comparing FSOLVE vs. FMINCON

①.1 Taylor-rule

Fmincon usually gets more solutions, even where fsolve says no sol.

①.2 RE-TC

Same

①.3 Anch-TC

Same.

I think this just means that fmincon is a little more easy-going w/ the sol. crit, since fsolve seeks to set $F(x) = 0$, while fmincon just tries to min $F(x)$

FSOLVE never finds a sol (it stalls once: $\{x, i\}_+$ in RE-TC, and init.)

FMINCON always finds at least 1 or 2 local mins.

But they are of 3 groups

- nonsensical: extremely volatile & huge
- really close to TR

↳ for TR & RE-TC → makes sense b/c you're either implementing the TR or the TR is a good implementation of RE-TC

- less volatile for x & i and more for π for anch-TC
↳ anch-TC calls for a less aggressive TR!

I actually think I should prioritize the "approximating the reaction function" approach of Peter over the value function iteration b/c it's more promising in terms of results.

But first I try reoptimizing by inputting the first $\{i_+^*\}$ as an initial guess for $\{i_t\}$.

→ doesn't work either.

command-approx-reaction.m

Problems:

✓ g_{π} is blowing up → a problem for the previous exercise too.

• Besides, in previous ex., was still evaluating the and TC wrong b/c at each t , need to account for CB's $E(\cdot)$ of future shocks

• loss isn't decreasing in any direction.

✓ forgot g_{π} .

$$\frac{\partial}{\partial \pi} \frac{1}{(dfe)^2} = \frac{\partial fe}{\partial \pi} \left(-2 \frac{1}{(dfe)^2} fe^{-1} \right)$$

$$= 2 \frac{1}{(dfe)^2} fe^{-1} \leftarrow =: g_{\pi}^{-1}$$

The problem w/ this is that $fe^{-3} \rightarrow \infty$ if fe is small.

I've inverted g_{π} and $g_{\bar{\pi}}$ b/c in the materials, I've defined $k_{+}^{-1} = g(\cdot)$, while in the code I have $k_{+} = g(\cdot)$.

Next: need to be smart about anchoring fe .

Correct evaluation of anchTC in the solve/fruition exercise. Regenerate.

$$fe = \pi_{+} - (\pi_{+-1} + b_{S_{+-1}})$$

$$k_{+} = g(fe) = \frac{1}{dfe^2}$$

$$k_{+}^{-1} = g(fe)^{-1} = dfe^2$$

$$\frac{\partial k_{+}^{-1}}{\partial \pi_{+}} = 2dfe =: g_{\pi}$$

$$\frac{\partial k_{+}^{-1}}{\partial \pi_{+-1}} = -2dfe =: g_{\bar{\pi}}$$

vs.

$$g_{\pi}^{-1}$$

vs.

$$g_{\bar{\pi}}^{-1}$$

- Still loss isn't decreasing

- Still TC eval is wrong for previous.

↳ I think I fixed that now - it takes way longer to run \rightarrow 25-30 min?

\rightarrow Took 17 minutes! (47 iter)
Or up to 30!

While it's running, let's work thru Colvard's VFI

$$u(c) = \frac{c^{1-\beta} - 1}{1-\beta}$$

$$k' = k^\alpha - c + (1-\delta)k$$

$$\Rightarrow V(k) = \max_c \frac{c^{1-\beta} - 1}{1-\beta} + \beta V(k'), \text{ or,}$$

plugging in $k'(k)$ for c ,

$$V(k) = \max_{k'} \frac{(k^\alpha + (1-\delta)k - k')^{1-\beta} - 1}{1-\beta} + \beta V(k')$$

Let's make sure that for no good value of $k(i)$, $k'(j)$

is $c < 0 \rightarrow$ for any $k(i)$, $k^\alpha(i) + (1-\delta)k(i) - k'(j) \stackrel{!}{\geq} 0$

I think I'm still evaluating the TC wrong 10 April 2020
- in all exercises, b/c if you input zero innovations,
that's gonna change f_a & f_b too. No - it's
correct b/c those are f_b & f_a that the CB expects
people to have. Phew!

Now I have the problem that `sim_learnLH.m`
doesn't work. It doesn't produce IRFs, and the
simulation doesn't seem to converge to RE either.

Sorry - it was b/c I turned σ & ρ of shocks to 0.
IRFs are working.

And sorry, it is converging, I plotted t/\bar{t}
and it does.

I'm cleaning up!

`sim_learnLH.m` is no longer touched!

`sim_learnLH_clean.m` = `sim_learnLH.m` and is meant

as a basis for subsequent work. (`materials24.m`

compares the two.)

sim-leamLL-deam-smooth.m is meant for
the smooth anchoring function only!

Ryan meeting

10 April 2020

→ should get a resid of zero if init at Tupper
rule sequence.

→ feed fsolve the T residuals and not the
square but the level

whereas fmincon will need the square.

Don't use fmincon b/c it takes lots of info out
lsqrnonlin since it uses a scalar

↳ similar to fmincon

Fsolve is the way to go when you know a
resid of 0 exists. Else lsqrnonlin b/c
it won't hit 0.

fsolve

1st thing to check: # of exog vars & equations
→ you can check that mistake in EE or
in the last period

A brute force: if $T=200$, $H=100$,
then at $t=150$, the EL in 2nd line of (B.1)
is fixed.

An implementable TC sounds interesting for a
discussion. It can also be an option for
this optimization. But it's not a concern
for the CB unless communicating w/ the
public is a concern.