Reproduction code package for "Decarboninzation Investment Strategies in an Uncertain Climate"

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The working paper version of our paper is: Bauer, A. M., F. McIsaac, S. Hallegatte. *How Delayed Learning about Climate Uncertainty Impacts Decarbonization Investment Strategies*. World Bank Policy Research Working Paper No. WPS10743, World Bank Group, Washington DC, 2024.

General package overview

This set of codes reproduces all of the figures and analysis carried out in *Decarbonization Investment Strategies* in an *Uncertain Climate*. This package uses Gurobi, a commercial nonlinear programming solver that is free for academics, but may not be free for everyone. (It is unclear to me if it's freely available for all researchers or just researchers at universities. I imagine Gurobi will handle this on a case-by-case basis, so just shoot their customer support staff an email and they can help.)

Each code is assigned a number corresponding to the figure it creates. Note an si before the script name indicates that figure is in the *Supplementary Information*. So code 01_xxx.py makes Figure 1 from the *main text*, which shows our calibration of the marginal abatement cost curves, while code si01_xxx.py makes the Figure 1 from the *Supplementary Information*. Here is the full table for both versions:

Figure Desired	Code to Run	Notes
Figure 1: Example abatement investment model and "strawman" model output.	01_simplified_simulations.sh	-
Figure 2: Calibrating marginal abatement costs	02_mac_calibration.sh	-
Figure 3: Effect of delayed learning on aggregate policy cost	03_effect_of_learning_low_lin	ear.sh
Figure 4: Effect of delayed learning on the temporal distribution of spending	04_temporal_redistribution_lo	w_linear.sh
Figure 5: Effect of delayed learning on sectoral allocation of abatement investment	05_sectoral_response.sh	-
Figure 6: Effect of delayed learning on the carbon price	06_carbon_price_response.sh	-
Figure SI 1: Effect of delayed learning on aggregate policy cost including direct air capture technologies	si01_dac_effect_of_learning.s	h -
Figure SI 2: Impact of delayed learning on sectoral allocation of abatement investment when direct air capture technologies are present	si02_dac_vs_no_dac_comp.sh	-

Figure Desired	Code to Run	Notes
Figure SI 3: Effect of delayed learning on aggregate policy cost,	si03_effect_of_learning_emis.sh	
growing emissions baseline		
Figure SI 4: Effect of delayed	si04_temporal_redistribution_	emis sh
learning on the temporal	bioi_semporar_rearburibation_	CMID: DI
distribution of spending, growing		
emissions baseline		
Figure SI 5: Effect of delayed	si05_effect_of_learning_high_	linear.sh
learning on aggregate policy cost,		
high-bound calibration		
Figure SI 6: Effect of delayed	si06_temporal_redistribution_	high_linear.sh
learning on the temporal		
distribution of spending,		
high-bound calibration		
Figure SI 7: Effect of delayed	si07_effect_of_learning_pow.s	Sh This figure was verified virtually.
learning on aggregate policy cost, nonlinear calibration		See <i>Known issues</i> below.
Figure SI 8: Effect of delayed	si08_temporal_redistribution_	nou sh
learning on the temporal	S100_temporal_redistribution_	.pow.sii
distribution of spending,		
nonlinear calibration		
Figure SI 9: Effect of delayed	si09_effect_of_learning_t15.s	sh -
learning on aggregate policy cost,		
$T^* = 1.5 \deg C$		
Figure SI 10: Effect of delayed	si10_temporal_redistribution_	t15.sh
learning on the temporal		
distribution of spending, $T^*=1.5$		
deg C		
Figure SI 11: Carbon prices as a	sill_carbon_price_sensitivity	
function of learning date		See Known issues below.

If you're an academic, you can email Gurobi customer support to get a free academic license. It's easy to install, and once it's installed, I believe you'll be good to go to run the codes.

A final note is that you should consider using the .yml file provided in this directory to establish a virtual python environment that should include all of the necessary dependencies for the code to run smoothly. I recommend using conda or mamba to do this.

How to run the code

To run the codes, simply navigate to the codes directory and run the numbered code to recreate the desired figure. If you want to run the program script_name, you may need to execute:

```
chmod +x script_name
```

to grant execution permissions (hence the +x) to the script you want to run.

As an example, if you want to recreate Figure 1 which shows our calibration of the marginal abatement cost curves, you would simply run:

./02_mac_calibration.sh

Notice the first bit of the above program name, 02_mac_calibration.sh, matches the figure number we wanted to create, Figure 1.

All figures will be deposited into the codes/figs folder. To run individual simulations, you can run any of the files in simulation_mains, and to make individual figures, you can run any file in the figure_mains folder. Note: You should run all scripts from the codes directory. As an example, let's say you want to run the invBase_cvxpy_main.py file in the ar6_15 calibration, but not save the output. Then in your command line, you'd use:

python simulation mains/invBase cvxpy main.py ar6 15 1 0

Note: You should be operating in the Python environment provided at the head directory. Without it, I make no guarantees any of this will run on your machine (and even then, well, mileage may vary...).

Known issues

The only figures that were not able to be reproduced on a member of the World Bank Group's Reproducibility team's computer were scripts si07 and si11. si07 was verified virtually, with a team member joining via video call and watching the code run on the author's laptop. si11 remains slightly different between the reproducer's run and the lead authors.

The hypothesized reason is that the simulations required for si07 and si11 have highly convex objective functions, which requires more powerful hardware to solve precisely than what was available to the reproducibility team member. The team member got an optimal_inaccurate solution during optimization step of si07. The code will throw an error when anything other than an optimal solution is found. It was verified that the original author of the code gets an optimal solution when the si07 code is run.

For si11, what is plotted is the optimal carbon price for each run, which is the Lagrange dual of emissions reductions in the model. This is found by solving the KCP dual optimization problem, which is even more sensitive to highly nonlinear objective functions than the minimization step itself (which is required for si07 and throws errors for the World Bank team member). While the results for a given user may vary, the overall qualitative implications of the figure as discussed in the paper hold for both the World Bank team member and the original author.

If a user gets an optimal_inaccurate solution for either of these codes, one possible course of action is to edit the scale parameter found in the codes/simulation_mains/invRec_RiskPrem_cvxpy_main.py file. This can be found on lines 34 through 42. This issue of scaling the objective function is common in optimization.

The hardware of the original author is a 2023 MacBook Pro with an M2 Pro Chip and 16 GB of RAM. Last edited: 8 May, 2025.