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Generate 500 observations from an AR(1) process Y_t with $\mathbb{E}[Y_t] = 0$, $\phi = 0.4$ and the variance of the white-noise forcing term $\sigma_{\varepsilon}^2 = 0.2$ using the two methods below (hint: the random number generator for i.i.d. normal is randn).

a. A for loop using the recursive structure of the AR(1).

Solution.

We study the stationary AR(1) process

$$Y_t = \mu + \phi (Y_{t-1} - \mu) + \varepsilon_t, \qquad \varepsilon_t \sim \mathcal{N}(0, \sigma^2), \quad |\phi| < 1.$$
 (1)

In the experiment, we set T=500, $\phi=0.4$, $\sigma^2=0.2$, $\mu=0$, and $Y_0=0$. A single sequence of shocks $\{\varepsilon_t\}_{t=1}^T$ is generated once and reused across implementations, enabling a pathwise comparison.

```
1 rng(12345, 'twister');
2 T
         = 500;
         = 0.4;
3 phi1
4 \text{ sigma2}_1 = 0.2;
         = 0;
                      % E[Y_t] = 0
5 mu1
                      % matching starting condition for both methods
6 YO
         = 0;
8 % Use the SAME innovation sequence for both methods
9 eps1 = sqrt(sigma2_1) * randn(T,1);
11 % (a) For-loop simulation
12 Y_loop = simulate_ar1_loop(T, phi1, sigma2_1, mu1, Y0, eps1);
13
14 %% From the "Helper functions" section:
function Y = simulate_ar1_loop(T, phi, sigma2, mu, Y0, eps)
17 %SIMULATE_AR1_LOOP Simulate AR(1) using explicit recursion (for-loop).
     Y_t = mu + phi*(Y_{t-1} - mu) + eps_t
18 %
19 %
      Inputs:
          T, phi, sigma2, mu, YO -> scalars
20 %
21 %
          eps -> T-by-1 vector of innovations (optional)
22 %
      Output:
23 %
          Y
               -> T-by-1 simulated series
24
      if nargin < 6 || isempty(eps)</pre>
25
           eps = sqrt(sigma2) * randn(T,1);
      end
26
               = zeros(T,1);
27
      Y(1)
               = mu + phi*(Y0 - mu) + eps(1);
28
      for t = 2:T
29
          Y(t) = mu + phi*(Y(t-1) - mu) + eps(t);
30
31
      end
32 end
```

The function simulate_ar1_loop implements the law of motion directly:

$$Y_1 = \mu + \phi (Y_0 - \mu) + \varepsilon_1, \qquad Y_t = \mu + \phi (Y_{t-1} - \mu) + \varepsilon_t \ (t = 2, \dots, T).$$

Passing the precomputed innovation vector ensures that any differences with alternative implementations are not driven by different random draws.

b. Using the function filter.

Solution.

Let $X_t := Y_t - \mu$. Then

$$(1 - \phi L) X_t = \varepsilon_t \iff X_t = \phi X_{t-1} + \varepsilon_t,$$

with L the lag operator. MATLAB's filter returns the zero-state solution Z_t to $(1 - \phi L)Z_t = \varepsilon_t$ (i.e., it implicitly sets $X_0 = 0$). To match an arbitrary initial condition $X_0 = Y_0 - \mu$, we add the homogeneous component:

$$X_t = Z_t + \phi^t X_0,$$

and then recover $Y_t = X_t + \mu$. The function simulate_ar1_filter implements precisely this decomposition so, under the same $\{\varepsilon_t\}$ and Y_0 , it reproduces the loop path exactly, period by period.

```
Y_filt = simulate_ar1_filter(T, phi1, sigma2_1, mu1, Y0, eps1);
3 %% From the "Helper functions" section:
5 function Y = simulate_ar1_filter(T, phi, sigma2, mu, Y0, eps)
6 %SIMULATE_AR1_FILTER Simulate AR(1) using FILTER for the centered process.
      Let X_t = Y_t - mu, then X_t = phi*X_{t-1} + eps_t.
      We generate X via filter and then shift back by mu.
     This implementation adds the exact initial-condition term phi^t * X0
9 %
10 %
      so the result matches the loop simulation pointwise.
      if nargin < 6 || isempty(eps)
11
          eps = sqrt(sigma2) * randn(T,1);
12
      end
13
      XO = YO - mu;
                           % initial condition for centered process
14
      % Zero-initial-condition filtered component
15
      Z = filter(1, [1, -phi], eps);
16
      t = (1:T)';
      X = Z + (phi.^t) * X0;  % exact IC adjustment
18
      Y = X + mu;
19
20 end
```

c. Check that when the forcing variables are the same, the output of the two approaches is the same (be careful with the starting conditions and with the random number generator). Solution.

To verify equivalence we compute

$$\max_{1 \le t \le T} |Y_t^{\text{loop}} - Y_t^{\text{filter}}|,$$

which is numerically zero up to machine precision.

```
diff_vec = Y_loop - Y_filt;
max_abs_diff = max(abs(diff_vec));
fprintf('[Ex.1] Max |difference| between methods: %.3e\n', max_abs_diff);

% Plot: overlay the two series
fh1 = figure('Position',[100 100 800 400]);
plot(1:T, Y_loop, '-', 'DisplayName','Loop'); hold on
plot(1:T, Y_filt, '--', 'DisplayName','Filter'); grid on
xlabel('t'); ylabel('$Y_t$')
title('Exercise 1: AR(1) via Loop vs. Filter (Overlay)')
legend('Location','best')
```

```
12 exportFig(fh1,'ex1_overlay.png');
13
14 % Plot: difference
15 fh2 = figure('Position',[100 100 800 350]);
16 plot(1:T, diff_vec, '-'); grid on
17 xlabel('t'); ylabel('$Y^{loop}_t - Y^{filter}_t$')
18 title(sprintf('Exercise 1: Difference, Max = %.1e', max_abs_diff))
19 exportFig(fh2,'ex1_difference.png');
```

We also provide two diagnostic figures: an overlay of the two series in Figure 1-1a and the path of their difference in Figure 1-1b.

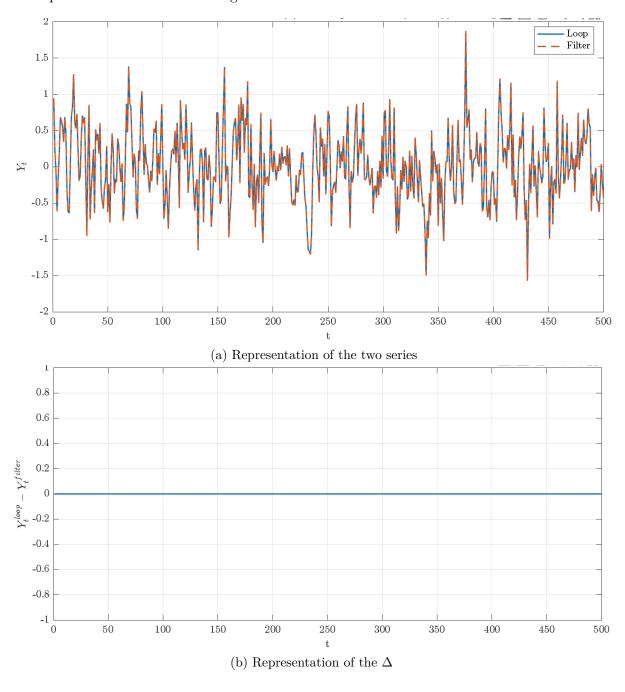


Figure 1-1: Visual representations supporting the statement that the two series are equivalent

Generate data from an AR(1) with $\phi = 0.6$, $\sigma_{\varepsilon}^2 = 0.4$ and $\mathbb{E}[Y_t] = 3$. Set the starting condition of your simulation to 20. What happens if the starting condition you choose is far from the unconditional mean of the process? What would you do in order to make sure that the sample path is a "proper" realization of the stationary process you want to simulate from? You can use either for or filter.

Solution.

We consider the AR(1)

$$Y_t = \mu + \phi (Y_{t-1} - \mu) + \varepsilon_t, \qquad \varepsilon_t \sim \mathcal{N}(0, \sigma^2), \quad |\phi| < 1, \tag{2}$$

with parameters $\phi = 0.6$, $\sigma^2 = 0.4$, $\mu = 3$. The horizon is T = 500 and the initial condition is set far from the mean, $Y_0 = 20$.

Writing $X_t := Y_t - \mu$ yields $X_t = \phi X_{t-1} + \varepsilon_t$, so the exact solution is

$$X_t = \phi^t X_0 + \sum_{j=0}^{t-1} \phi^j \varepsilon_{t-j}, \quad \text{hence} \quad \mathbb{E}[Y_t \mid Y_0] = \mu + \phi^t (Y_0 - \mu).$$
 (3)

With $\phi = 0.6$ and $Y_0 - \mu = 17$, the deterministic component $\phi^t(Y_0 - \mu)$ decays geometrically: the half-life is $h = \log(1/2)/\log(\phi) \approx 1.36$ periods, so the expected path returns rapidly toward $\mu = 3$. Figure 2-2a shows a simulated trajectory together with the unconditional mean.

```
% AR(1) with phi=0.6, sigma^2=0.4, E[Y_t]=3, start Y0=20
3 rng(23456, 'twister');
         = 500;
5 phi2
          = 0.6;
6 \text{ sigma2}_2 = 0.4;
7 \text{ mu2} = 3;
                       % unconditional mean
                       % starting far from mean
8 \text{ YO}_{\text{far}} = 20;
10 % Simulate with a for-loop (explicit control over initial condition)
11 eps2 = sqrt(sigma2_2) * randn(T,1);
      = simulate_ar1_loop(T, phi2, sigma2_2, mu2, Y0_far, eps2);
14 % Plot the sample path and the unconditional mean
15 fh3 = figure('Position',[100 100 900 360]);
plot(1:T, Y2, '-', 'DisplayName','$Y_t$'); hold on; grid on
yline(mu2, '--', '$\mathrm{E}[Y_t]=\mu=3$', 'Interpreter','latex', '
      LabelVerticalAlignment','bottom', 'DisplayName', '$\mathrm{E}[Y_t]=\mu=3$')
xlabel('t'); ylabel('$Y_t$')
19 title ('Exercise 2: AR(1) Path with Initial Condition Far from Mean')
20 legend('Location','best')
21 exportFig(fh3,'ex2_path_far_from_mean.png');
```

To obtain a draw that is effectively free of initial-condition transients, we simulate T+B periods and discard the first B (burn-in). Since $|\phi| < 1$, the effect of Y_0 on Y_t is proportional to ϕ^t ; choosing B = 500 makes ϕ^B negligible. The retained segment $\{Y_{B+1}, \ldots, Y_{B+T}\}$ is therefore well-approximated by a sample from the stationary distribution. Figure 2-2b displays the post-burn-in path with the mean line.

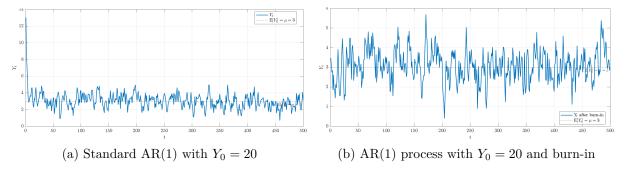


Figure 2-2: Visual representations supporting the statement that the two series are equivalent

Compute the empirical distribution of the OLS estimator in the case of an AR(1) with $\phi = 0.4$ and T = 250 (you are free to choose the variance of the innovation). Construct a t-test for the null hypothesis $H_0: \phi = 0$, against a two-sided alternative $H_1: \phi \neq 0$. How often do you reject H_0 at the 95% confidence level when T = 250?

Solution.

We consider the AR(1) data-generating process (DGP)

$$Y_t = \phi Y_{t-1} + \varepsilon_t, \qquad \varepsilon_t \sim \mathcal{N}(0, \sigma^2),$$
 (4)

with $\phi = 0.4$, $\sigma^2 = 1$, and T = 250. To remove transients from the initial condition, each replication simulates T + B observations with a burn-in of B = 300 and retains the last T points. We run R = 5000 replications.

```
% Empirical distribution of OLS estimator; t-test of HO: phi=0
  % DGP: AR(1) with phi=0.4, T=250.
4 rng(34567, 'twister');
         = 250;
         = 0.4;
  sigma2_3 = 1.0;
                        % explicit
         = 0;
  mu3
         = 5000;
                        % number of Monte Carlo replications
9 R
                        % short burn-in for stationarity
         = 300;
10
phi_hat = zeros(R,1);
13 tstat
        = zeros(R,1);
```

¹With $|\phi| < 1$, the effect of Y_0 on Y_t decays like ϕ^t ; a burn-in of 300 makes this negligible.

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In each replication we estimate ϕ by OLS from the regression $Y_t = \phi Y_{t-1} + u_t$ (no constant since the DGP has mean zero):

$$\widehat{\phi} = \frac{\sum_{t=2}^{T} Y_{t-1} Y_t}{\sum_{t=2}^{T} Y_{t-1}^2}, \qquad \widehat{u}_t = Y_t - \widehat{\phi} Y_{t-1}.$$

Let $\hat{s}^2 = \sum_{t=2}^T \hat{u}_t^2/(T-1)$ and $\operatorname{se}(\hat{\phi}) = \sqrt{\hat{s}^2/\sum_{t=2}^T Y_{t-1}^2}$. We test $H_0: \phi = 0$ with the usual t-statistic

$$t = \frac{\widehat{\phi} - 0}{\operatorname{se}(\widehat{\phi})},$$

and reject for $|t| > t_{0.975,T-1}$ (two-sided 5%).

```
for r = 1:R
      % Innovations and simulation length with burn-in
      TT = T + B;
      eps3 = sqrt(sigma2_3) * randn(TT,1);
      % Start at the mean (mu3) + burn-in
6
7
      Ytmp = simulate_ar1_loop(TT, phi3, sigma2_3, mu3, mu3, eps3);
           = Ytmp(B+1:end);
                                              % keep last T observations
8
9
      \% OLS in Y_t = phi * Y_{t-1} + u_t (no intercept; mean is zero)
      ylag = Y(1:end-1);
           = Y(2:end);
12
          = ylag;
                                               % (T-1) \times 1
13
      bhat = (X' * X) \ (X' * yt);
14
      uhat = yt - X * bhat;
16
      % --- Correct degrees of freedom: nu = (T-1) - 1 = T - 2 ---
17
      nu = (T - 1) - 1;
18
           = (uhat' * uhat) / nu;
                                               % unbiased sigma_u^2
19
           = sqrt( s2 / (X' * X) );
                                               % std error of bhat
20
21
      phi_hat(r) = bhat;
      tstat(r) = bhat / se;
                                               % test H0: phi = 0
23
```

The Monte Carlo summary reports the empirical mean and standard deviation of $\hat{\phi}$ across replications and the rejection frequency at the 5% level. In our run, the mean of $\hat{\phi}$ is slightly below the true value (a familiar small finite-sample downward bias for positive ϕ), while the 5% test of H_0 : $\phi = 0$ rejects essentially always, reflecting very high power at T = 250 and $\phi = 0.4$.

Remark 1 As $T \to \infty$, $\sqrt{T}(\hat{\phi} - \phi) \stackrel{d}{\to} \mathcal{N}(0, 1 - \phi^2)$, so, under $H_1: \phi \neq 0$, the t-statistic is approximately normal with noncentrality

$$\lambda \approx \frac{\phi}{\sqrt{(1-\phi^2)/T}} = \phi \sqrt{\frac{T}{1-\phi^2}}.$$

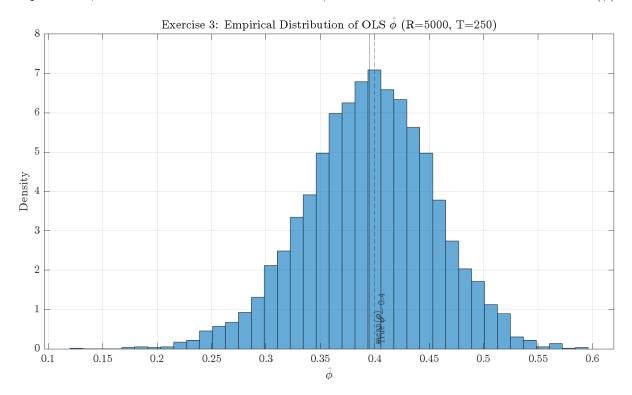
For $\phi = 0.4$ and T = 250, this gives $\lambda \approx 6.9$, implying near–unit power, consistent with the simulated rejection rate.

```
1 % Rejection frequency at 5%
2 if exist('tinv','file')
3    tcrit = tinv(0.975, nu);
4 else
5    tcrit = 1.96; % approximation for moderate T
6 end
7 reject = mean(abs(tstat) > tcrit);
```

Mean $(\hat{\phi})$	0.3949
Standard Deviation $(\hat{\phi})$	0.0589
reject_H0_rate	1

Table 3-1: Summary statistics

Figure 3-3: The histogram below depicts the empirical sampling distribution of $\widehat{\phi}$ across the R replications, with vertical lines at the true value $\phi = 0.4$ and at the Monte Carlo mean mean $(\widehat{\phi})$.



Compute the empirical distribution of the OLS estimator in the case of an AR(1) with $\phi = 0.9$ and $T \in \{50, 100, 200, 1000\}$. For each T, do 1000 simulations and plot the distribution. How is the distribution changing with T?

Solution.

For each sample size $T \in \{50, 100, 200, 1000\}$ we simulate R = 1000 samples, discarding a burn–in of B = 500 observations to remove dependence on initial conditions. For each replication we compute the OLS estimator from the regression through the origin (consistent with $\mu = 0$):

$$\widehat{\phi} = \frac{\sum_{t=2}^{T} y_{t-1} y_t}{\sum_{t=2}^{T} y_{t-1}^2}.$$

We also record the usual OLS standard error and t-statistic for testing $H_0: \phi = 0$ (formally using n = T - 1 observations and $\nu = n - 1 = T - 2$ degrees of freedom).

In our notes, the least-squares probability limit for an AR(1) is

$$plim \hat{\phi} = \frac{\gamma_1}{\gamma_0} = \phi,$$

where $\gamma_h = \text{Cov}(y_t, y_{t-h})$. Thus $\widehat{\phi}$ is consistent. Moreover, under standard regularity conditions,

$$\sqrt{T}(\widehat{\phi} - \phi) \stackrel{d}{\to} \mathcal{N}(0, 1 - \phi^2),$$

so that, asymptotically,

$$\operatorname{sd}(\widehat{\phi}) \approx \sqrt{\frac{1-\phi^2}{T}}.$$

For $\phi = 0.9$, $1 - \phi^2 = 0.19$, so $sd(\hat{\phi}) \approx \sqrt{0.19/T}$.

The Monte Carlo follows exactly this design:

- 1. Draw $\{\varepsilon_t\}_{t=1}^{T+B}$ i.i.d. $\mathcal{N}(0,1)$.
- 2. Generate $\{y_t\}$ recursively with the given ϕ , keep the last T observations after the burn-in.
- 3. Compute $\hat{\phi}$ via the no-intercept OLS formula above.
- 4. For inference quantities, we use n=T-1 effective observations and an unbiased residual variance estimator $\hat{\sigma}^2 = \text{RSS}/(n-1) = \text{RSS}/(T-2)$, which implies $\text{se}(\hat{\phi}) = \sqrt{\hat{\sigma}^2/\sum y_{t-1}^2}$ and t-critical values based on $\nu = T-2$ degrees of freedom.

```
14
15 for iT = 1:numel(Ts)
      T = Ts(iT);
16
      nu = T - 2;
                                         % df for regression with (T-1) rows, 1
17
      slope
18
      phi_hat = zeros(R,1);
19
      tstat = zeros(R,1);
20
21
      for r = 1:R
22
          % --- simulate AR(1) with burn-in
23
          TT = T + B;
24
          eps = sqrt(sigma2_4) * randn(TT,1);
25
          Yall = simulate_ar1_loop(TT, phi, sigma2_4, mu, mu, eps);
26
27
                = Yall(B+1:end);
                                         % keep last T observations
          \% --- OLS: Y_t = phi * Y_{t-1} + u_t (no intercept since mu=0)
          ylag = Y(1:end-1);
               = Y(2:end);
31
          уt
               = ylag;
                                        % (T-1)-by-1 regressor
32
          Х
          XX = X' * X;
33
34
          bhat = XX \setminus (X' * yt);
35
          uhat = yt - X * bhat;
36
37
          % --- correct finite-sample variance and t-stat
38
                                      % unbiased residual variance: RSS/(T-2)
          s2 = (uhat' * uhat) / nu;
          se = sqrt( s2 / XX );
                                          % std error of slope
41
          tstat(r)
                    = bhat / se;
                                        % test H0: phi = 0
42
          phi_hat(r) = bhat;
43
      end
44
      \% --- Monte Carlo summaries for this T
45
          = mean(phi_hat);
46
           = std(phi_hat);
47
      sd
      bias = m - phi;
48
49
      if exist('tinv','file')
                                          % two-sided 5% test against HO: phi = 0
          tcrit = tinv(0.975, nu);
51
52
          tcrit = 1.96;
                                          % normal approx
53
54
      rej = mean(abs(tstat) > tcrit);
55
56
      E4_summary{iT,:} = [T, m, sd, bias, rej];
57
58
      % --- Histogram for this T
      fh = figure('Position',[100 100 840 420]);
      histogram(phi_hat, 40, 'Normalization','pdf'); hold on; grid on
61
      xline(phi, '--', 'True $\phi=0.9$', 'LabelVerticalAlignment','bottom', '
62
      Interpreter','latex');
      xline(m, '-', '$\mathrm{mean}(\hat{\phi})$', 'LabelVerticalAlignment','
63
      bottom', 'Interpreter', 'latex');
      xlabel('$\hat{\phi}$', 'Interpreter','latex');
64
      ylabel('Density', 'Interpreter','latex');
65
      title(sprintf('Exercise 4: OLS on AR(1), $\\phi=0.9$ (T=%d, R=%d)', T, R),
66
      'Interpreter', 'latex');
      exportFig(fh, sprintf('ex4_hist_T%d.png', T));
68 end
69
70 % Save table
71 writetable(E4_summary, fullfile(outdir,'ex4_summary.csv'));
```

From the R = 1000 replications we obtain, for each T, the Monte Carlo mean, standard

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deviation, and bias (mean minus 0.9). The reported summary is:

T	$\operatorname{mean}(\widehat{\phi})$	$\operatorname{sd}(\widehat{\phi})$	bias	$\operatorname{rej}\{H_0\colon \phi=0\}$
50	0.8682	0.0756	-0.0318	1.000
100	0.8840	0.0493	-0.0160	1.000
200	0.8908	0.0322	-0.0092	1.000
1000	0.8983	0.0138	-0.0017	1.000

Two key comparisons with the asymptotic variance formula help interpret these numbers. Theoretical $sd(\phi) \approx \sqrt{0.19/T}$ yields 0.062 (T=50), 0.044 (T=100), 0.031 (T=200), and 0.0138 (T=1000). The Monte Carlo standard deviations are close and converge toward these values as T grows (the mild over-dispersion at small T is a known finite-sample feature). The Monte Carlo mean shows the familiar downward small-sample bias for $\phi > 0$, which is O(1/T) and vanishes as T increases.

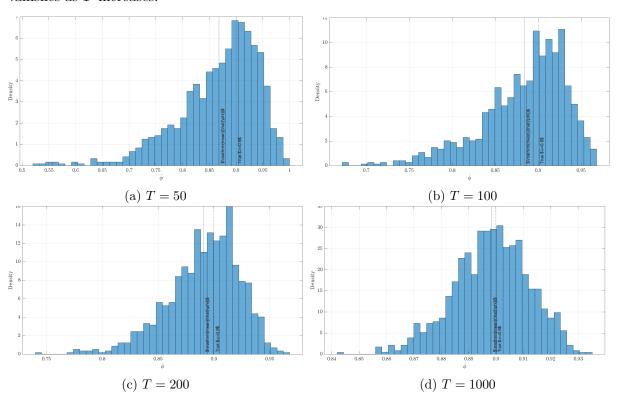


Figure 4-4: Evolution of the distributions as the number of observations T increases.

The histograms (one per T) and the table convey the same message:

- Centering: The distribution is centered below ϕ for small T (negative finite-sample bias), but the bias shrinks rapidly (from about -0.032 at T = 50 to about -0.002 at T = 1000), in line with consistency.
- Spread: The dispersion decreases roughly at the $\propto T^{-1/2}$ rate. Numerically, the standard deviation falls from $\approx 0.076~(T=50)$ to $\approx 0.014~(T=1000)$, very close to $\sqrt{(1-\phi^2)/T}$.
- Shape: By the central limit reasoning above, the standardized estimator $\sqrt{T}(\hat{\phi} \phi)$ becomes approximately normal as T increases. Empirically, the histograms become more symmetric and concentrated around 0.9 as T grows, with thinner tails.

Compute the empirical distribution (do 1000 simulations) of the OLS estimator in the regression $x_t = ax_{t-1} + v_t$ in the case in which the data-generating process for x_t is MA(1) with $\theta = 0.6$ for $T \in \{50, 100, 200, 1000\}$. What is the mean of the distributions? Does the mean converge to anything as $T \to \infty$? Discuss.

Solution.

We simulate MA(1) data

$$x_t = \varepsilon_t + \theta \varepsilon_{t-1}, \qquad \varepsilon_t \stackrel{i.i.d.}{\sim} \mathcal{N}(0, \sigma^2), \ \theta = 0.6,$$

keep the last T observations after a burn-in of B = 500, and estimate, for each replication,

$$\hat{a} = \arg\min_{a} \sum_{t=2}^{T} (x_t - a x_{t-1})^2 = \frac{\sum_{t=2}^{T} x_{t-1} x_t}{\sum_{t=2}^{T} x_{t-1}^2}.$$

We repeat this R = 1000 times for each T, plot the histogram of \hat{a} , and compute the Monte Carlo mean and standard deviation.

Because the regression omits the MA structure, OLS converges to the *projection* coefficient of x_t on x_{t-1} :

plim
$$\hat{a} = \frac{\text{Cov}(x_t, x_{t-1})}{\text{Var}(x_{t-1})} = \frac{\gamma_1}{\gamma_0} = \rho(1).$$

For an MA(1), $\gamma_0 = (1 + \theta^2)\sigma^2$ and $\gamma_1 = \theta\sigma^2$, so

$$p\lim \hat{a} = \frac{\theta}{1 + \theta^2}.$$

With $\theta = 0.6$, this gives

$$\frac{0.6}{1+0.6^2} = \frac{0.6}{1.36} = 0.441176\overline{47}.$$

```
_1 % OLS of x_t on x_{t-1} when x_t is MA(1) with theta=0.6
2 \% DGP: x_t = eps_t + theta * eps_{t-1}, eps_t ~ N(0, sigma^2)
4 rng(56789, 'twister');
                                        % Reproducibility for Ex.5
5 \text{ theta} = 0.6;
6 \text{ sigma2}_5 = 1.0;
                          % explicit
      = [50, 100, 200, 1000];
       = 1000;
       = 500;
                           % burn-in for MA(1)
9 B
11 % Theoretical plim of OLS when regressing x_t on x_{t-1}: rho(1) = theta/(1+
     theta<sup>2</sup>)
plim_a = theta / (1 + theta^2);
14 E5_summary = table('Size',[numel(Ts) 5], ...
       'VariableTypes',{'double','double','double','double','double'}, ...
16
      'VariableNames', {'T', 'mean_a_hat', 'sd_a_hat', 'bias_from_plim', '
      theoretical_plim', });
17
  for iT = 1:numel(Ts)
      T = Ts(iT);
19
      a_hat = zeros(R,1);
20
21
      for r = 1:R
22
           TT = T + B;
23
           % Simulate MA(1) with burn-in
```

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Stefano Graziosi
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```
[x_all, ~] = simulate_ma1(TT, theta, sigma2_5, 0);
25
26
          x = x_all(B+1:end);
          Xlag = x(1:end-1); xt = x(2:end);
          bhat = (Xlag' * Xlag) \ (Xlag' * xt); % no constant
          a_hat(r) = bhat;
      end
31
      m = mean(a_hat);
33
      sd = std(a_hat);
34
      bias = m - plim_a;
35
      E5_summary{iT,:} = [T, m, sd, bias, plim_a];
36
37
      % Plot histogram for this T
38
      fh = figure('Position',[100 100 840 420]);
      histogram(a_hat, 40, 'Normalization','pdf'); hold on; grid on
      xline(plim_a, '--', sprintf('plim = %.3f', plim_a), 'LabelVerticalAlignment
      ','bottom');
                     '-', '$\mathrm{mean}(\hat{a})$', 'LabelVerticalAlignment','
      xline(m,
42
      bottom');
      xlabel('$\hat{a}$'); ylabel('Density')
43
      title(['Exercise 5: OLS on MA(1) Data, ', '$\theta=0.6$', ' (T=', num2str(T
44
      ), ', R=', num2str(R), ')'], 'Interpreter', 'latex')
      exportFig(fh, sprintf('ex5_hist_T%d.png', T));
45
46 end
47
48 % Save table
49 writetable(E5_summary, fullfile(outdir,'ex5_summary.csv'));
51 %% From the "Helper functions" section:
53 function [x, eps] = simulate_ma1(T, theta, sigma2, mu)
54 %SIMULATE_MA1 Simulate MA(1): x_t = mu + eps_t + theta*eps_{t-1}
      if nargin < 4, mu = 0; end</pre>
      eps = sqrt(sigma2) * randn(T,1);
56
      % Vectorized MA(1): set eps_0 = 0 and use lagged innovations
57
      eps_lag = [0; eps(1:end-1)];
59
      x = mu + eps + theta * eps_lag;
60 end
```

Across the R=1000 replications, the empirical results are as follows:

T	$\operatorname{mean}(\hat{a})$	$\operatorname{sd}(\hat{a})$	Bias from plin	Theoretical plim
50	0.434493054710814	0.106511187264365	-0.00668341587742133	0.441176470588235
100	0.437911528963784	0.0742479286576639	-0.00326494162445146	0.441176470588235
200	0.435441067223447	0.0535341394861661	-0.00573540336478806	0.441176470588235
1000	0.439680841416116	0.0230883947393155	-0.00149562917211887	0.441176470588235

Table 5-2: Summary statistics for different sample sizes T.

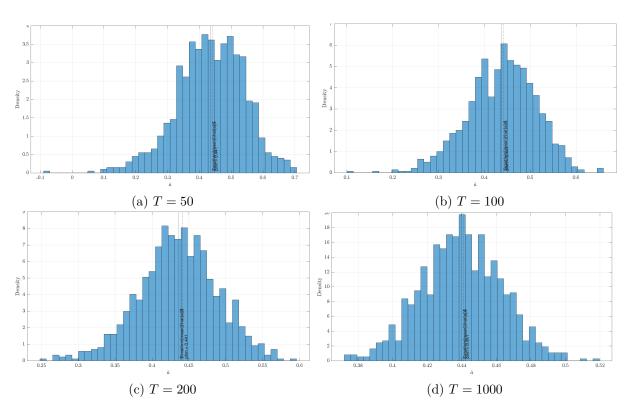


Figure 5-5: Evolution of the distributions as the number of observations T increases.

Focusing on the mean, we have:

$$\begin{array}{c|ccccc} T & 50 & 100 & 200 & 1000 \\ \hline mean(\hat{a}) & 0.4345 & 0.4379 & 0.4354 & 0.4397 \end{array}$$

with empirical standard deviations approximately (0.1065, 0.0742, 0.0535, 0.0231), respectively.

In light of these results, we can go back to the original question and discuss what happens to the mean(\hat{a}):

- Center (mean). The Monte Carlo mean of \hat{a} is close to 0.4412 and drifts toward it as T increases. Thus, the mean converges to the pseudo-true parameter $\rho(1) = \theta/(1 + \theta^2)$, not to a structural AR(1) coefficient (there is no AR(1) here).
- Spread (variance). The dispersion of \hat{a} shrinks with T at the usual $\propto T^{-1/2}$ rate (histograms become more concentrated and more nearly normal around 0.4412).
- Interpretation. Regressing x_t on x_{t-1} is estimating the best linear predictor slope at lag 1. For MA(1) data, this equals the lag-1 autocorrelation. Hence, in large samples the OLS fit recovers $\rho(1)$.

```
1 % Optional: plot mean(\hat{a}) across T for a compact summary figure
2 fh = figure('Position',[100 100 620 360]); grid on; hold on
3 plot(E5_summary.T, E5_summary.mean_a_hat, '-o', 'DisplayName','$\mathrm{mean}(\hat{a})$');
4 yline(plim_a, '--', 'DisplayName','theoretical plim');
5 xlabel('$T$'); ylabel('Mean of $\hat{a}$ across replications')
6 title('Exercise 5: Convergence of $\mathrm{mean}(\hat{a})$ with $T$', '
Interpreter','latex')
7 legend('Location','best')
8 exportFig(fh,'ex5_mean_vs_T.png');
```

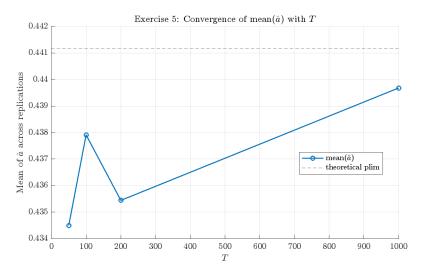


Figure 5-6: Evolution of the mean(\hat{a})

Write a function that generates T observations from an ARMA(p,q) using the for-loop approach. The function must have the following inputs:

- (i) the number of observations;
- (ii) the variance of the white noise ε_t
- (iii) the coefficients of the AR and MA polynomials or the roots of the AR and MA polynomials (hint: you may find the poly and roots functions useful);

and as output the realizations of the ARMA(p,q) and the white noise ε_t . Solution.

We tried to include all of the necessary informations and explanations directly in the function description down below.

```
function [y, e] = arma_sim(T, sigma2, ar_in, ma_in, varargin) %#ok<DEFNU>
2
  %ARMA_SIM Generate T observations from an ARMA(p,q) via for-loop.
3
      y_t = mu + sum_{i=1}^p phi_i (y_{t-i} - mu) + e_t + sum_{j=1}^q theta_j e_{t-i}
     t-j}
4
  %
      with e_t \sim N(0, sigma2).
5
6
  %
    Inputs (required):
7
  %
               -> number of observations to RETURN (after optional burn-in)
  %
               -> variance of white noise e_t
8
      sigma2
               -> AR parameters: either coefficients [phi_1..phi_p] OR lag-roots
  %
      ar in
9
      [lambda_1..lambda_p]
10 %
      ma in
               -> MA parameters: either coefficients [theta_1..theta_q] OR lag-
     roots [lambda_1..lambda_q]
11 %
12 % Name-Value pairs (optional):
13 %
      'ParamType' -> 'coeffs' (default) or 'roots'.
14 %
                     If 'roots', we interpret:
                       A(L) = prod_{i=1}^p (1 - lambda_i L) => 1 - phi_1 L - ...
15 %
      - phi_p L^p
                       B(L) = prod_{j=1}^q (1 + lambda_j L) => 1 + theta_1 L +
16 %
      \dots + theta_q L^q
17 %
                     Coefficients are then recovered from these lag polynomials.
18 %
      'BurnIn'
                  -> number of burn-in observations to discard (default 500)
19 %
                  -> unconditional mean mu (default 0)
```

```
20 %
21 % Outputs:
      y -> T-by-1 vector of ARMA(p,q) observations
      e -> T-by-1 vector of shocks e_t used to generate y
25 % Notes:
    * Stationarity (AR) / invertibility (MA) are the user's responsibility.
26 %
      * We simulate with burn-in (default 500) from zero initial conditions.
27 %
28
      p = numel(ar_in); q = numel(ma_in);
29
      ip = inputParser; ip.KeepUnmatched = true;
30
      addParameter(ip,'ParamType','coeffs');
31
      addParameter(ip,'BurnIn',500);
32
33
      addParameter(ip,'Mu',0);
      parse(ip, varargin{:});
      paramType = validatestring(ip.Results.ParamType, {'coeffs','roots'});
      B = ip.Results.BurnIn; mu = ip.Results.Mu;
37
38
      % Determine phi and theta
39
      switch paramType
40
          case 'coeffs'
41
               phi = ar_in(:).';
                                              % row
42
               theta = ma_in(:).';
43
           case 'roots'
44
               % Build lag polynomials and read off implied coefficients.
45
               if p>0
                   poly_ar = 1; % A(L)
47
48
                   for i=1:p
                       poly_ar = conv(poly_ar, [1, -ar_in(i)]); % (1 - lambda_i L)
49
50
                   phi = -poly_ar(2:end);  % A(L) = 1 - phi_1 L - ... - phi_p L^p
               else
                   phi = [];
54
               if q>0
                   poly_ma = 1; \% B(L)
                   for j=1:q
                       poly_ma = conv(poly_ma, [1, ma_in(j)]); % (1 + lambda_j L)
58
59
                   theta = poly_ma(2:end); % B(L) = 1 + theta_1 L + ... + theta_q
      L^q
               else
61
                   theta = [];
62
63
      end
64
      % Sanity: warn if (approx) nonstationary / noninvertible
      if ~isempty(phi)
67
          A = [1, -phi(:).'];
68
          rr = roots(A);
69
          if any(abs(rr) <= 1)</pre>
70
               warning('AR polynomial has roots at or inside unit circle; process
      may be nonstationary.');
72
          end
73
      if ~isempty(theta)
74
          Bpoly = [1, theta(:).'];
          rr = roots(Bpoly);
           if any(abs(rr) <= 1)
               warning('MA polynomial has roots at or inside unit circle; process
78
      may be noninvertible.');
79
          end
```

```
end
80
81
       % Simulation with burn-in
82
       TT = T + B; p = numel(phi); q = numel(theta);
83
       e = sqrt(sigma2) * randn(TT,1);
84
       y = zeros(TT, 1);
       if mu ~= 0
86
           % Work with deviations from mu for numerical stability
87
           x = zeros(TT,1); % x_t = y_t - mu
88
           for t = 1:TT
89
               accAR = 0; accMA = 0;
90
               for i=1:p
91
                    if t-i \ge 1, accAR = accAR + phi(i) * x(t-i); end
92
93
94
               for j=1:q
                    if t-j \ge 1, accMA = accMA + theta(j) * e(t-j); end
               x(t) = accAR + e(t) + accMA;
97
           end
98
           y = x + mu;
99
       else
100
           for t = 1:TT
101
               accAR = 0; accMA = 0;
102
               for i=1:p
103
                    if t-i \ge 1, accAR = accAR + phi(i) * y(t-i); end
104
105
               for j=1:q
                    if t-j \ge 1, accMA = accMA + theta(j) * e(t-j); end
107
108
               y(t) = accAR + e(t) + accMA;
109
           end
       end
111
112
       % Drop burn-in
113
       y = y(B+1:end);
114
       e = e(B+1:end);
115
116 end
```