```
a) Cointegration
clear ; close all; clc
load Data Canada
Y = Data(:, 3:end);
figure
plot(dates, Y, 'LineWidth', 2)
xlabel('Year')
ylabel('Percent')
names = series(3:end);
legend(names, 'location', 'NW')
title('{\bf Canadian Interest Rates, 1954-1994}')
axis tight
grid on
y1 = Y(:,1); % Short-term rate
% Levels data:
fprintf('=== Test y1 for a unit root ===\n\n')
[h1,pVal1] = adftest(y1,'model','ARD') % Left-tail probability
fprintf('\n=== Test y1 for stationarity ===\n\n')
[h0,pVal0] = kpsstest(y1,'trend',false) % Right-tail probability
% Differenced data:
fprintf('\n=== Test (1-L)y1 for a unit root ===\n\n')
[h1D,pVal1D] = adftest(diff(y1), 'model', 'ARD') % Left-tail
probability
fprintf('\n=== Test (1-L)y1 for stationarity ===\n\n')
[h0D,pVal0D] = kpsstest(diff(y1),'trend',false) % Right-tail
probability
figure
plot(dates(2:end), diff(Y), 'LineWidth', 2)
names = series(3:end);
legend(names, 'location', 'NW')
title('{\bf Differenced Data}')
axis tight
grid on
% Run the test with both "tau" (t1) and "z" (t2) statistics:
```

```
fprintf('\n=== Engle-Granger tests for cointegration ===\n\n')
[hEG,pValEG] = egcitest(Y,'test',{'t1','t2'})
% Return the results of the cointegrating regression:
[\sim, \sim, \sim, \sim, reg] = egcitest(Y, 'test', 't2');
c0 = reg.coeff(1);
b = reg.coeff(2:3);
figure
C = get(gca, 'ColorOrder');
set(gca,'NextPlot','ReplaceChildren','ColorOrder',circshift(C,3));
plot(dates, Y*[1;-b]-c0, 'LineWidth', 2);
title('{\bf Cointegrating Relation}');
axis tight
grid on
% Permutations of the data variables:
P0 = perms([1 2 3]);
[~,idx] = unique(P0(:,1)); % Rows of P0 with unique regressand y1
P = P0(idx,:); % Unique regressions
numPerms = size(P,1);
% Preallocate:
T0 = size(Y, 1);
HEG = zeros(1, numPerms);
PValEG = zeros(1, numPerms);
CIR = zeros(T0, numPerms);
% Run all tests:
for i = 1:numPerms
    YPerm = Y(:, P(i,:));
    [h,pVal,~,~,reg] = egcitest(YPerm,'test','t2');
    HEG(i) = h;
    PValEG(i) = pVal;
    c0i = reg.coeff(1);
    bi = reg.coeff(2:3);
    CIR(:,i) = YPerm*[1;-bi]-c0i;
end
fprintf('\n=== Different Engle-Granger tests, same data ===\n\n')
HEG, PValEG
```

```
% Plot the cointegrating relations:
figure
C = get(gca, 'ColorOrder');
set(qca,'NextPlot','ReplaceChildren','ColorOrder',circshift(C,3))
plot(dates,CIR,'LineWidth',2)
title('{\bf Multiple Cointegrating Relations}')
legend(strcat({'Cointegrating relation '}, ...
     num2str((1:numPerms)')),'location','NW');
axis tight
grid on
fprintf('\n=== Johansen tests for cointegration ===\n')
[hJ,pValJ] = jcitest(Y,'model','H1','lags',1:2);
[~,~,~,~,mles] = jcitest(Y,'model','H1','lags',2,'display','params');
B = mles.r2.paramVals.B % Cointegrating relations with rank = 2
restriction
fprintf('\n=== Test y1, y2, y3 for stationarity ===\n\n')
[h0J,pVal0J] = jcontest(Y,1,'BVec',{[1 0 0]',[0 1 0]',[0 0 1]'})
clear;
echo off;
N = 500;
                             % Number of particles.
T = 100;
                             % Number of time steps.
                            % Continuous state dimension.
n x = 4;
                            % Number of discrete states.
n z = 3;
                            % Dimension of observations.
n y = 4;
parameter. A = zeros(n x, n x, n z);
parameter.B = zeros(n x, n x, n z);
parameter.C = zeros(n y, n x, n z);
parameter.D = zeros(n_y, n_y, n_z);
parameter.E = zeros(n x, n x, n z);
parameter.F = zeros(n x, 1, n z);
parameter.G = zeros(n y, 1, n z);
ST = 1;
                       % sampling period
u = ones(1,T);
                                         % input signal(Deterministic)
```

```
b) particle filter, rao-blackwellized particle filter, extended kalman
%Defining parameters
for i=1:n z
  parameter.A(:,:,i) = [1 ST 0 0; 0 1 0 0; 0 0 1 ST; 0 0 0 1];
  parameter.C(:,:,i) = eye(n x);
  parameter.B(:,:,i) = 0.2*eye(n x,n x);
  parameter.D(:,:,i) = sqrt(3)*diag([20,1,20,1]);
 parameter.G(:,:,i) = zeros(n y,1);
% input and control vectors as given in paper
parameter.F(:,1,1) = [0 \ 0 \ 0]';
parameter.F(:,1,2) = [-1.225, -0.35, 1.225, 0.35]';
parameter.F(:,1,3) = [1.225, 0.35, -1.225, -0.35]';
% Transition matrix
parameter.T = [0.9 \ 0.05 \ 0.05;
                0.05 0.9 0.05;
                0.05 0.05 0.91;
                                                     % Initial r
parameter.pz0 = [0 1 0];
markov chain
parameter.mu0 = zeros(n x, 1);
                                               % Initial Gaussian mean
                                               % Initial Gaussian
parameter.S0 = 1*eye(n x, n x);
covariance
% Generating ground truth
x = zeros(n x, T);
z = zeros(1,T);
y = zeros(n y,T);
x(:,1) = parameter.mu0 + sqrtm(parameter.S0)*randn(n x,1);
z(1) = length(find(cumsum(parameter.pz0')<rand))+1;</pre>
for t=2:T
  z(t) = length(find(cumsum(parameter.T(z(t-1),:)') < rand)) + 1;
  x(:,t) = parameter.A(:,:,z(t))*x(:,t-1) +
parameter.B(:,:,z(t))*randn(n x,1) + parameter.F(:,:,z(t))*u(:,t);
  y(:,t) = parameter.C(:,:,z(t))*x(:,t) +
parameter.D(:,:,z(t))*randn(n y,1) + parameter.G(:,:,z(t))*u(:,t);
end
```

```
%%%Plotting ground truth and noisy observations
figure; hold on
plot(x(1,:), x(3,:), '*');
plot(y(1,:), y(3,:), '.');
drawnow
[styles, colors, symbols, str] = plotColors;
%%%% Plotting change in the state of target as it maneuvers
figure; hold on
for k=1:3
  ndx=find(z==k);
 plot(x(1,ndx), x(3,ndx), sprintf('%s%s', colors(k), symbols(k)));
  plot(x(1,ndx), x(3,ndx));
end
title('data')
drawnow
%%%% particle filter jmls
[zest pf, xest pf, zamples pf, xsamples pf] = pfSlds(N, par, y, u);
%%%% Rao-Balckwellized particle filter
[zest rbpf, xest rbpf, zsamples rbpf] = rbpfSlds(N, par, y, u);
%%%% Extended Kalman filter
[zest ekf, xest ekf, zsamples ekf] = ekf(N, par, y, u);
%%%%Calculating Mean square error
mse ekf = mean((xest ekf(1,:)-x(1,:)).^2) +
mean((xest ekf(3,:)-x(3,:)).^2);
mse pf = mean((xest pf(1,:)-x(1,:)).^2) +
mean((xest pf(3,:)-x(3,:)).^2);
mse rbpf = mean((xest rbpf(1,:)-x(1,:)).^2) +
mean((xest rbpf(3,:)-x(3,:)).^2);
[junk, z max pf] = max(zest pf, [], 1);
%%%%Plotting extended Kalman filter estimate
[junk,z max ekf] = max(zest ekf,[],1);
figure; hold on
for k=1:n z
 ndx=find(z max ekf==k);
```

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plot(xest ekf(1,ndx), xest ekf(3,ndx));
end
title(sprintf('ekf, mse %5.3f', mse ekf));
%%%Plotting Particle filter estimate
figure; hold on
for k=1:n z
 ndx=find(z max pf==k);
 plot(xest pf(1,ndx), xest pf(3,ndx));
end
title(sprintf('pf, mse %5.3f', mse pf));
%%%%Plotting rbpf estimate
[junk,z max rbpf] = max(zest rbpf,[],1);
figure; hold on
for k=1:n z
  ndx=find(z max rbpf==k);
 plot(xest rbpf(1,ndx), xest rbpf(3,ndx));
end
title(sprintf('rbpf, mse %5.3f', mse rbpf));
{
figure;
plot(1:T,z,'k',1:T,z,'ko',1:T,z max rbpf,'r+',1:T,z max pf,'bv','line
width',1);
legend('','True state','RBPF MAP estimate','PF MAP estimate');
% axis([0 T+1 0.5 n z+0.5])
% error rate Z
detect error pf = sum(z\sim=z max pf)/T;
detect error rbpf = sum(z\sim=z max rbpf)/T;
z ind = dummyEncoding(z(:), n z);
figure;
subplot(1,3,1); imagesc(z ind);
title('truth');
subplot(1,3,2); imagesc(zest pf');
title(sprintf('pf, error rate %5.3f', detect error pf));
subplot(1,3,3); imagesc(zest rbpf');
title(sprintf('rbpf, error rate %5.3f', detect error rbpf));
subplot(1,4,4); imagesc(zest ekf');
title(sprintf('ekf, error rate %5.3f', detect error ekf));
```

```
fprintf('PF: misclassification rate %5.3f, mse %5.3f, time %5.3f\n',
  detect error pf, mse pf, time pf);
fprintf('RBPF: misclassification rate %5.3f, mse %5.3f, time
%5.3f\n', ...
  detect error rbpf, mse rbpf, time rbpf);
fprintf('EKF: misclassification rate %5.3f, mse %5.3f, time %5.3f\n', ...
  detect error ekf, mse ekf, time ekf);
function [zest, xest, zsamples, w] = rbpfSlds(N, par, y, u,
resamplingScheme)
if nargin < 5, resamplingScheme = 2; end
 [nz,T] = size(y);
 [n y, n x, n z] = size(parameter.C);
 z rbpf = ones(1,T,N);
z rbpf pred = ones(1,T,N);
                                % One-step-ahead predicted values of
z.
mu = 0.01*randn(n x, T, N);
                                % Kalman mean of x.
mu pred = 0.01*randn(n x,N);
Sigma = zeros(n x, n x, N);
                                % Kalman covariance of x.
Sigma pred = zeros(n x, n x, N);
S = zeros(n y, n y, N);
                                 % Kalman predictive covariance.
y \text{ pred} = 0.01 \cdot \text{randn}(n y, T, N);
w = ones(T, N);
                                 % Importance weights.
initz = 1/n z*ones(1, n z);
xest = zeros(n x,T); % KPM
zest = zeros(n z,T); % KPM
for i=1:N
  Sigma(:,:,i) = 1*eye(n x, n x);
  Sigma pred(:,:,i) = Sigma(:,:,i);
  z rbpf(:,1,i) = length(find(cumsum(initz')<rand))+1;</pre>
  S(:,:,i) =
parameter.C(:,:,z) rbpf(1,1,i) *Sigma pred(:,:,i) *parameter.C(:,:,z) rb
pf(1,1,i))' + ...
parameter.D(:,:,z rbpf(1,1,i))*parameter.D(:,:,z rbpf(1,1,i))';
end
```

```
for t=2:T
  f(RBPF : t = \%i / \%i \ \ \ \ f('\n');
  % Sequential Importance Sampling Step:
  for i=1:N
    % sample z(t) \sim p(z(t) | z(t-1))
    z \text{ rbpf pred}(1,t,i) =
length(find(cumsum(parameter.T(z rbpf(1,t-1,i),:)')<rand))+1;</pre>
    % Kalman prediction:
    \operatorname{mu} \operatorname{pred}(:,i) = \operatorname{parameter.A}(:,:,z \operatorname{rbpf} \operatorname{pred}(1,t,i)) * \operatorname{mu}(:,t-1,i) +
. . .
                     parameter.F(:,:,z rbpf pred(1,t,i))*u(:,t);
Sigma pred(:,:,i)=parameter.A(:,:,z) rbpf pred(1,t,i))*Sigma(:,:,i)*pa
rameter.A(:,:,z) rbpf pred(1,t,i))'...
parameter.B(:,:,z rbpf pred(1,t,i))*parameter.B(:,:,z rbpf pred(1,t,i))
))';
    S(:,:,i) =
parameter.C(:,:,z rbpf pred(1,t,i))*Sigma pred(:,:,i)*parameter.C(:,:
,z rbpf pred(1,t,i))' + ...
parameter.D(:,:,z rbpf pred(1,t,i))*parameter.D(:,:,z rbpf pred(1,t,i)
))';
    y pred(:,t,i) = parameter.C(:,:,z rbpf pred(1,t,i))*mu pred(:,i)
                        parameter.G(:,:,z) rbpf pred(1,t,i)*u(:,t);
  end
  % Evaluate importance weights.
  for i=1:N
    w(t,i) = (det(S(:,:,i))^{(-0.5)})^* \dots
              \exp(-0.5*(y(:,t)-y_pred(:,t,i))'*pinv(S(:,:,i))*(y(:,t)-
. . .
                                      y \text{ pred}(:,t,i))) + 1e-99;
  end
% w(t,:) = \exp(\log w(t,:)) + 1e-99*ones(size(w(t,:)));
  w(t,:) = w(t,:)./sum(w(t,:)); % Normalise the weights.
  % Selection Step:
    if resamplingScheme == 1
    outIndex = residualR(1:N,w(t,:)');
```

```
elseif resamplingScheme == 2
    outIndex = deterministicR(1:N,w(t,:)');
  else
    outIndex = multinomialR(1:N,w(t,:)');
  z rbpf(1,t,:) = z rbpf pred(1,t,outIndex);
 mu pred = mu pred(:,outIndex);
  Sigma pred = Sigma pred(:,:,outIndex);
  S = S(:,:,outIndex);
  y pred(:,t,:) = y pred(:,t,outIndex);
  % Kalman Update:
    for i=1:N
    % Kalman update:
    K =
Sigma pred(:,:,i) *parameter.C(:,:,z rbpf(1,t,i)) '*pinv(S(:,:,i));
    mu(:,t,i) = mu pred(:,i) + K*(y(:,t)-y pred(:,t,i));
    Sigma(:,:,i) = Sigma pred(:,:,i) -
K*parameter.C(:,:,z rbpf(1,t,i))*Sigma pred(:,:,i);
  end
 xest(:,t) = mean(squeeze(mu(:,t,:)), 2); % KPM - unweighted mean
  zest(:,t) = normalize(hist(squeeze(z rbpf(1,t,:)), 1:n z)); % KPM
end % End of t loop.
zsamples = squeeze(z rbpf); % (1,t,:) \rightarrow (t,:)
end
function [zest, xest, zsamples, xsamples, w] = pfSlds(N, par, y, u,
resamplingScheme)
if nargin < 5, resamplingScheme = 2; end
 [nz,T] = size(y);
 [n y, n x, n z] = size(parameter.C);
z pf = ones(1,T,N);
                               z pf pred = ones(1,T,N);
One-step-ahead predicted values of z.
x pf = 10*randn(n x,T,N);
x pf pred = x pf;
y pred = 10*randn(n y,T,N);
w = ones(T, N);
                                % Importance weights.
```

```
initz = 1/n z*ones(1, n z);
xest = zeros(n x,T); %KPM
zest = zeros(n z,T); % KPM
for i=1:N
  z pf(:,1,i) = length(find(cumsum(initz')<rand))+1;</pre>
end
for t=2:T
  % Sequential Importance Sampling Step:
  for i=1:N
    % sample z(t) \sim p(z(t) | z(t-1))
    z pf pred(1,t,i) =
length(find(cumsum(parameter.T(z pf(1,t-1,i),:)')<rand))+1;</pre>
    % sample x(t) \sim p(x(t) | z(t|t-1), x(t-1))
   x 	ext{ pf pred}(:,t,i) = parameter.A(:,:,z 	ext{ pf pred}(1,t,i)) *
x pf(:,t-1,i) + ...
                       parameter.B(:,:,z pf pred(1,t,i))*randn(n x,1)
+ ...
                       parameter.F(:,:,z) pf pred(1,t,i)*u(:,t);
  end
  % Evaluate importance weights.
  for i=1:N
    y \text{ pred}(:,t,i) = \text{parameter.C}(:,:,z \text{ pf pred}(1,t,i)) *
x pf pred(:,t,i) + ...
                     parameter.G(:,:,z) pf pred(1,t,i)*u(:,t);
   Cov =
parameter.D(:,:,z pf pred(1,t,i))*parameter.D(:,:,z pf pred(1,t,i))';
   w(t,i) = (det(Cov)^{(-0.5)})*exp(-0.5*(y(:,t)-y pred(:,t,i))'* ...
                          pinv(Cov)*(y(:,t)-y pred(:,t,i))) + 1e-99;
  w(t,:) = w(t,:)./sum(w(t,:)); % Normalise the weights.
  % Selection Step:
  if resamplingScheme == 1
    outIndex = residualR(1:N,w(t,:)');
  elseif resamplingScheme == 2
```

```
outIndex = deterministicR(1:N,w(t,:)');
  else
    outIndex = multinomialR(1:N,w(t,:)');
  z pf(1,t,:) = z pf pred(1,t,outIndex);
  x pf(:,t,:) = x pf pred(:,t,outIndex);
  xest(:,t) = mean(squeeze(x pf(:,t,:)), 2); % KPM - unweighted mean
  zest(:,t) = normalize(hist(squeeze(z pf(1,t,:)), 1:n z)); % KPM
end % End of t loop.
zsamples = squeeze(z pf); % (1,t,:) \rightarrow (t,:)
xsamples = x pf;
end
function [X ind, X3d] = dummyEncoding(X, nStates)
[N, D] = size(X);
if nargin < 2
    nStates = nunique(X);
end
응 {
offset = cumsum(nStates);
offset = [0, offset(1:end-1)];
X = bsxfun(@plus, X, offset)';
I = repmat(1:N, D, 1);
K = max(sum(nStates), max(X(:)));
ndx = sub2ind([N, K], I(:), X(:));
X \text{ ind} = false(N, K);
X \text{ ind(ndx)} = \text{true;}
응 }
K = max(nStates);
X ind = zeros(N, sum(nStates));
X3d = zeros(N, D, K);
for d = 1:D
  idx = sum(nStates(1:d-1))+1:sum(nStates(1:d));
  miss = isnan(X(:,d));
  X ind(~miss,idx) = bsxfun(@eq, X(~miss,d), [1:nStates(d)]);
  X \text{ ind}(miss, idx) = NaN;
  X3d(~miss,d,1:nStates(d)) = reshape(X ind(~miss, idx), [sum(~miss)
1 nStates(d));
end
```

```
test = false;
if test
    X = bsxfun(@minus, X', offset); % return X back to original state
    if nargin < 2
         nStates = zeros(1,D);
         for j=1:D
             nStates(j) = length(unique(X(:,j)));
         end
    end
    X \text{ ind2} = zeros(N, sum(nStates));
    offset = 0;
    for s = 1:length(nStates)
         for i = 1:N
             X \text{ ind2}(i, \text{offset} + X(i, s)) = 1;
         end
         offset = offset+nStates(s);
    end
    toc
    assert(isequal(X ind, X ind2));
end
end
function [zest, xest, zsamples] = ekf(N, par, y, u)
[nz,T] = size(y);
[n y, n x, n z] = size(par.C);
 z = kf = ones(1,T,N);
z = kf pred = ones(1,T,N);
mu = 0.01*randn(n x, T, N);
mu pred = 0.01*randn(n x,N);
Sigma = zeros(n x, n x, N);
Sigma pred = zeros(n x, n x, N);
S = zeros(n y, n y, N);
y \text{ pred} = 0.01 \times \text{randn}(n y, T, N);
w = ones(T, N);
initz = 1/n z*ones(1, n z);
xest = zeros(n x,T); % KPM
zest = zeros(n z,T); % KPM
for i=1:N
  Sigma(:,:,i) = 1*eye(n x, n x);
  Sigma pred(:,:,i) = Sigma(:,:,i);
  z ekf(:,1,i) = length(find(cumsum(initz')<rand))+1;</pre>
```

```
S(:,:,i) =
par.C(:,:,z = kf(1,1,i))*Sigma pred(:,:,i)*par.C(:,:,z = kf(1,1,i))' + ...
                                       par.D(:,:,z ekf(1,1,i))*par.D(:,:,z ekf(1,1,i))';
end
          for i=1:N
          % sample z(t) \sim p(z(t) | z(t-1))
          z = kf(1,t,i) = length(find(cumsum(par.T(z = kf(1,t-1,i),:)') < rand))+1;
% Kalman prediction:
         mu pred(:,i) = par.A(:,:,z ekf pred(1,t,i))*mu(:,t-1,i) + ...
                                              par.F(:,:,z ekf pred(1,t,i))*u(:,t);
Sigma pred(:,:,i) = par.A(:,:,z ekf pred(1,t,i))*Sigma(:,:,i)*par.A(:,:,z ekf pred(1,t,i))*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:,:,i)*Sigma(:
f pred(1,t,i))'...
par.B(:,:,z ekf pred(1,t,i))*par.B(:,:,z ekf pred(1,t,i))';
         S(:,:,i) =
par.C(:,:,z ekf pred(1,t,i))*Sigma pred(:,:,i)*par.C(:,:,z ekf pred(1,t,i)
) ' + ...
                                  par.D(:,:,z ekf pred(1,t,i))*par.D(:,:,z ekf pred(1,t,i))';
         y \text{ pred}(:,t,i) = par.C(:,:,z \text{ ekf } pred(1,t,i))*mu \text{ pred}(:,i) + ...
                                                     par.G(:,:,z ekf pred(1,t,i))*u(:,t);
    % Evaluate importance weights.
    for i=1:N,
         w(t,i) = (det(S(:,:,i))^{(-0.5)})^* \dots
                               \exp(-0.5*(y(:,t)-y \text{ pred}(:,t,i))**pinv(S(:,:,i))*(y(:,t)-...
                                                                                        y \text{ pred}(:,t,i))) + 1e-99;
    end;
% w(t,:) = \exp(\log w(t,:)) + 1e-99*ones(size(w(t,:)));
    w(t,:) = w(t,:)./sum(w(t,:)); % Normalise the weights.
       for i=1:N.
         % Kalman update:
         K = Sigma pred(:,:,i)*par.C(:,:,z ekf(1,t,i))'*pinv(S(:,:,i));
         mu(:,t,i) = mu pred(:,i) + K*(y(:,t)-y pred(:,t,i));
         Sigma(:,:,i) = Sigma pred(:,:,i) -
K*par.C(:,:,z ekf(1,t,i))*Sigma pred(:,:,i);
    xest(:,t) = mean(squeeze(mu(:,t,:)), 2); % KPM - unweighted mean
    zest(:,t) = normalize(hist(squeeze(z ekf(1,t,:)), 1:n z)); % KPM
end; % End of t loop.
```

```
zsamples = squeeze(z ekf); % (1,t,:) \rightarrow (t,:)
end
function [styles, colors, symbols, str] = plotColors()
% Use plot(x,y,str{i}) to print in i'th style
% Use colors and linestyles, not markers
colors = ['b' 'r' 'k' 'q' 'c' 'y' 'm' ...
        'r' 'b' 'k' 'g' 'c' 'y' 'm'];
symbols = ['o' 'x' '*' '>' '<' '^' 'v' ...
        '+' 'p' 'h' 's' 'd' 'o' 'x'];
styles = {'-', ':', '-.', '--', ':', '-.', '--', ...
  '-', ':', '-.', '--', ':', '-.', '--'};
for i=1:length(colors)
  str\{i\} = sprintf('-%s%s', colors(i), symbols(i));
  str{i} = sprintf('sss', colors(i), styles{i});
end
end
function outIndex = deterministicR(inIndex,q)
if nargin < 2, error('Not enough input arguments.'); end
[S,arb] = size(q); % S = Number of particles.
% Residual Resampling:
N off= zeros(1,S);
u=zeros(1,S);
cumDist = cumsum(q');
aux=rand(1);
u=aux:1:(S-1+aux);
u=u./S;
j=1;
for i=1:S
   while (u(1,i)>cumDist(1,j))
      j=j+1;
   end
   N off(1,\dot{j})=N off(1,\dot{j})+1;
end
```

```
%figure;
%plot(cumDist,'-b'); hold on;
%plot(u, '-r');
%figure;
%bar(N off);
index=1;
for i=1:S
  if (N \circ ff(1,i) > 0)
    for j=index:index+N off(1,i)-1
      outIndex(j) = inIndex(i);
    end
  end
  index= index+N off(1,i);
end
function outIndex = multinomialR(inIndex,q)
if nargin < 2, error('Not enough input arguments.'); end
[S,arb] = size(q); % S = Number of particles.
% Multinomial Sampling:
N off= zeros(1,S);
cumDist= cumsum(q');
% generate S ordered random variables uniformly distributed in [0,1]
u = fliplr(cumprod(rand(1,S).^(1./(S:-1:1))));
\dot{j} = 1;
for i=1:S
  while (u(1,i)>cumDist(1,j))
    j = j + 1;
  N off(1,j)=N off(1,j)+1;
end
index=1;
for i=1:S
  if (N \circ ff(1,i) > 0)
    for j=index:index+N_off(1,i)-1
      outIndex(j) = inIndex(i);
```

```
end
  end
  index= index+N off(1,i);
end
end
function outIndex = residualR(inIndex,q)
    if nargin < 2, error('Not enough input arguments.'); end
[S,arb] = size(q); % S = Number of particles.
% vanilla re-sample:
N off= zeros(1,S);
% first integer part
q res = S.*q'; %'
N off = fix(q res);
% residual number of particles to sample
N res=S-sum(N off);
if (N res \sim = 0)
 q res=(q res-N off)/N res;
  cumDist= cumsum(q res);
  % generate N res ordered random variables uniformly distributed in
[0,1]
  u = fliplr(cumprod(rand(1, N res).^(1./(N res:-1:1))));
  j=1;
  for i=1:N res
    while (u(1,i)>cumDist(1,j))
      j=j+1;
    N off(1,j)=N off(1,j)+1;
  end
end
%copy traj
index=1;
for i=1:S
  if (N \circ ff(1,i) > 0)
    for j=index:index+N off(1,i)-1
      outIndex(j) = inIndex(i);
    end
  end
  index= index+N off(1,i);
end
```

(Particle filter has been implemented close to Dunham which is then extended to rbpf and $\ensuremath{\mathsf{ekf}}\xspace$)