Probabilistic Inference and Learning to Control: pilco

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Abstract

The pilco toolbox is an Octave and Matlab implementation of time series inference and controller optimisation for nonlinear dynamical systems in continuous-state, discrete-time settings. The toolbox is both flexible and extensible. Flexibility is achieved by allowing user specification of which dynamics models, inference procedures and policy functional forms to use. Users may choose to provide their own data of a dynamical system, or generate synthetic data using any of the accompanying 'scenarios' such as the cart-pole system. Extensibility is a result of the modular code structure, separating class types such as controllers, dynamics models, scenarios, each of which are easily added to.

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1 Overview

There are several supporting structures, classes and functions the user should be aware of. Each is discussed in further details in later sections. Many functions will have derivative counterparts, named similarly except with an additional suffix 'd'.

1.1 Toolbox Components

Scenario: Scenarios are computer simulators of dynamical systems. Each scenario represents a particular dynamical system - such as the cart-pole system - and generates synthetic data using the known laws of motion. To begin a simulation, execute the **doit.m** file within a scenario's directory.

Dynamics Model: A dynamics model object is a probabilistic model which is used to make inferences from observed system evolutions and predictions on future dynamics. Each model is either a Gaussian process (GP), or GP state space model (GP-SSM).

Policy: A policy is a deterministic function, parameterised by θ , that maps state distribution s to action u; $\pi: s \times \theta \to u$. Various functional forms of the policy are available, including linear, trigonometric, Gaussian radial basis functions (RBFs) and combinations thereof.

Controller: A controller is a wrapper object of the policy function. In addition, a controller computes various covariances, derivatives and can filter the sequence of system-state observations.

Cost: The cost function takes a state distribution s and returns a scalar value on the unit interval; $c: s \to [0,1]$ The output represents the instantaneous cost (or measure of *undesirability*) per unit of time of the system being in state s. The cost specifies what it is the user wants to achieve, a 0 cost being the most desirable kind of state, and a 1 being the least desired.

1.2 Program Flow

Select policy function form, parameterised by (initially random) θ :

$$\pi: s \times \theta \to u.$$
 (1)

FOR each trial n in [1, 2, ..., N]:

- Infer dynamics given all observed data so far; $p(s^{t+1}|s^t, u^t, D^{1:n-1})$.
- Simulate system from initial state distribution s_0 until final state distribution $s_{horizon}$, using the dynamics model to make successive predictions from one time step to the next.
- Evaluate controller:

$$f(s_0; \theta) = \sum_{t=0}^{horizon} \gamma^t cost(s_t),$$
 (2)

where f is the cumulative discounted cost, θ is the policy parameters. Derivative information $\frac{df}{d\theta}$ is also computed.

• Optimise policy (via gradient decent):

$$\theta^* \quad \leftarrow \quad \underset{\theta}{\arg\min} \ f(s_0;\theta). \tag{3}$$

• Apply controller and generate more training data \mathcal{D}^n . This may be done using an available scenario, or the user's hardware.

1.3 States

1.3.1 State Representation

The *state* of the control system can be expressed in various ways. The simplest representation is the *physical state*, consisting:

- 1. present position vector $\mathbf{x}^{\mathbf{t}}$,
- 2. present velocity vector $\dot{\mathbf{x}}^{t}$.

For the best results, we incorporate all relevant information such that satisfies the Markov assumption. One way the above Markov assumption is violated is with laggy sensors. Under sensor lag, observations reflect the state of the system in the past, during which time the previous action has still been in effect. Thus previous actions become relevant when inferring the current state. So a better state representation when mitigating the effects of sensor-lag would be:

- 1. previous action vector \mathbf{u}^{t-1} ,
- 2. present position vector $\mathbf{x}^{\mathbf{t}}$,
- 3. present velocity vector $\dot{\mathbf{x}}^{t}$.

The above representation is especially important when sensor lag amounts to a significant proportion of the time discretisation. Now consider the case where, in addition to sensor delay, velocity information $\dot{\mathbf{x}}$ is unavailable. In this case, the 2-Markov state representation is helpful, consisting:

- 1. previous-previous action vector \mathbf{u}^{t-2} ,
- 2. previous position vector \mathbf{x}^{t-1} ,
- 3. previous action vector \mathbf{u}^{t-1} ,
- 4. present position vector $\mathbf{x}^{\mathbf{t}}$.

Intuitively, the above state representation works well because one could estimate any relevant velocity information using finite differences. All aforementioned state representations are possible in the pilco toolbox, including general N-Markov representations. Note the ordering of state variables must be chronological.

The *state distribution* is approximated as a multivariate Gaussian over all state variables. We encode the state distribution with a structure:

```
state % state structure
state.m % state mean vector
state.s % state covariance matrix
```

1.3.2 State Variable Ordering

Each scenario directory contains its own executable doit.m file, used to begin simulations. The doit file specifies the *ordering* of the state variables (amongst other things). For example, consider a hypothetical scenario: a 2-Markov cart-pole with an augmented variable. The state variables would be ordered as follows:

```
1 1 1 oou even older value of u
2 2 2 ox old cart position
3 3 otheta old angle of the pendulum
```

```
4
       4
                         old value of u
 4
           ou
    5
 5
       5
                         cart position
           Х
 6
       6
                         angle of the pendulum
           theta
 7
           v
                         cart velocity
 8
                         angular velocity
           dtheta
 9
           x - ox
                         relative change in cart position
10
                         force applied to cart
           u
       7
                         sine of old angle of the pendulum
           sin(otheta)
       8
           cos(otheta)
                         cosine of old angle of the pendulum
       9
                         sine of angle of the pendulum
           sin(theta)
                         cosine of angle of the pendulum
      10
           cos(theta)
```

where the green column of numbers are the indices for each state variable in our state (here our state dimensionality is D=6). The three types of indices always agree the first D numbers. The red column keeps extending past index D, and includes some extra variables that the simulator will use. The blue column, used by the policy, performs some trigonometric operations on some of the state variables, and thus also has an extends past D.

The red and green indices must refer to variables in a chronological order. The chronological order also takes the understanding that action variables at any time t occur *after* state variables at time t. E.g. index 10 is the final red index.

Typically a **doit** file will only display the red and green indices. The blue indices are specificed implicitly as a function of the green indices and the set of angular variables discussed below.

1.3.3 State Variable Classes

Let us now group the above variables that we used in the above subsection. There exists three different ways of indexing each state variable in Sec. 1.3.2 (red green and blue) and the sets of indices below have been colour coded w.r.t. which columns of indices they refer to.

```
D
                       % dimensionality of the state
     = 6;
Ε
    = 2;
                       % number of outputs from the dynamics model
    = 1;
                       % dimensionality of the action
angi = [3 6];
                           % indices for variables treated as angles (sin/cos)
augi = [9]:
                           % indices for variables augmented to ode vars
dyni = [1 2 3 4 5 6];
                           % indices for input into the dynamics model
dyno = [5 6];
                           % indices of dynamics model output, and loss input
odei = [5 6 7 8];
                           % indices for the ode solver
poli = [1 2 4 5 7 8 9 10]; % indices for the inputs to the policy
```

The dyni and dyno variables are currently redundant, since it is always the case that dyni = 1:D, and dyno = D-E+1:D.

Note the [7 8 9 10] indices of **poli** refer to the trigonometric terms of whatever **angi** specified were angle variables.

2 Base functions

The base directory contains common files that are always used to train a controller. Evaluation of a controller's parameterisation is with loss.m (Sec. 2.1). The controller's parameterisation loss is a function of the cumulative cost distribution, provided by simulate.m (Sec. 2.2). An important

subroutine of **simulate.m** is the **propagate.m** function (Sec. 2.3). The **propagate.m** function computes predictive state distributions at time step t + 1 given the state distribution at time t current time.

2.1 Loss

5a

```
\langle loss.m 5a \rangle \equiv
 1 function [f, df] = loss(p, s, dyn, ctrl, cost, H, exp, cc_prev, n)
2 % If the heuristic function 'explore' is not present (or if it is empty) then
3 % the expected cumulative cost is returned. If 'explore' is present, then the
4 % loss is a function of the cumulative cost mean and variance of the current
5 % parameterisation 'cc', and the parameterisation of the previous rollout
6 % 'cc_prev', and the number of rollout trials remaining 'n'. Derivatives of
7 % these quantities are computed when desired.
9 % [f, df] = loss(p, s, dyn, ctrl, cost, H, exp, cc_prev, n)
10 %
11 % p
                  policy parameter structure
12 % s
                  initial state structure
13 % dyn
                 dynamics model object
14 % ctrl
                  controlloer object
15 % cost
                  cost object
16 % H
                  time-steps horizon
17 % exp
                  exploration struct
18 % cc_prev cumulative (discounted) cost structure of previous rollout
                  number of trials remaining (inc. current point in time, so n>0)
19 % n
20 % f
            1x1 loss
21 % df
            1xP loss derivative wrt policy parameters
22 %
23 % Copyright (C) 2015 by Carl Edward Rasmussen and Rowan McAllister 2015-06-01
25 ctrl.set_policy_p(p);
26 exploring = exist('explore','var') && ~isempty(explore);
27 sargs = {s, dyn, ctrl, cost, H};
29 if nargout == 1 % no derivatives
    [S, A, ~, cc] = simulate(sargs{:});
30
31
    if exploring
      f = explore(exp, cc_prev, n, S, A, cc, [], sargs{:});
32
33
    else
34
      f = cc.m;
35
    end
36 else % derivatives
    [S, A, ~, cc, dcc] = simulate(sargs{:});
37
38
    if exploring
39
      [f, df] = explore(exp, cc_prev, n, S, A, cc, dcc, sargs{:});
40
   else
41
      f = cc.m;
42
      df = dcc.m;
43
    end
44 end
```

2.2 Simulate

The *cumulative cost* of a distribution over states is defined to be the long term (possibly discounted) total cost of starting in a given state distribution and following a policy π up to a time horizon

H. The simulate function returns states struct arrays, action struct arrays, cost struct arrays, the cumulative costs and possibly cumulative cost derivatives wrt the policy parameters.

```
\langle simulate.m 5b \rangle \equiv
5b
       1 function [s, a, c, cc, dcc] = simulate(s, dyn, ctrl, cost, H)
       3 % [s, a, c, cc, dcc] = simulate(s, dyn, ctrl, cost, H)
       4 %
       5 % s
                        . state structure F*1 mean vector
       6 % m
       7 % s
                          F*F covariance matrix
       8 % ?
                          possibly other fields representing additional information
      8 % ? possibly other freeds replaced and a dynamics model object

10 % ctrl controller object

11 % is struct indexing vectorized state distributions variables

12 % m F*1 indices of mean parameters

13 % s F*F indices of variance parameters

13 % s pumber of parameters in the policy
      14~\% np number of parameters in the policy 15~\% ns number of state distribution parameters.
                                  number of state distribution parameters (means and variances)
      16~\% policy . policy structure
      17 % p . policy parameters structure
18 % cost . cost function object
19 % cov @ function returning cost-covariances between two states
20 % fcn @ function returning expected and variance of a states's cost
      \frac{21\ \%}{22\ \%\ H} discount factor length of prediction horizon \frac{20\ \%}{23\ \%} s (output) H+1 state struct array containing Gaussian state distributions
      24~\% m $\mathrm{F*1}$ mean vector
      26 % a H action struct array containing Gaussian action distributions
27 % m U*1 mean vector
28 % s U*U covariance matrix
29 % c H+1 cost struct array containing Gaussian cost distributions
30 % m mean scalar
      31 % s
                                  variance scalar
      32 % cc
                                cumulative (discounted) cost structure
      33 % m
                                mean scalar
                         variance scalar
derivative structure of cc
      34 % s
      35 % dcc
      36 % m
                          1xP derivative cc-mean wrt policy parameters, same fields as p
      37 % s 1xP derivative cc-variance wrt policy parameters, same fields as p
      38 %
      39 % Copyright (C) 2008-2015 Carl Edward Rasmussen, 2015-05-31
      41 global currT; currT=1;
      42 if ~isfield(s,'s'); s.s = zeros(length(s.m)); end
      43 c = cost.fcn(s);
                                                                                                           % init c
      44 \text{ cc.m} = \text{c.m}; \text{ cc.s} = \text{c.s};
                                                                                                           % init cc
                                                                                                           % init q
      45 q = s.s;
      46 gamma = cost.gamma; D = ctrl.D; F = length(s.m);
      47 if nargout < 5 % no derivatives
      48 for t = 1:H
                                                                                    % iterate up to horizon
      49
             \langle compute \ state \ action \ cost \ 7a \rangle
               if nargout > 3 % then accumulate discounted cost
      50
                  ⟨compute cumulative costs 7b⟩
      51
      52
               end
      53 end
      54 else % do derivatives
```

```
55 \( \compute cumulative costs and derivatives 7c \)
56 \( \text{end} \)
```

The states, actions and cost distributions are computed by iterating up to the horizon, propagating forward the state distributions and recording the action distributions and computing the cost.

```
7a \langle compute \ state \ action \ cost \ 7a \rangle \equiv (5b)

1 [s(t+1), C, a(t)] = propagate(s(t), dyn, ctrl); % propagate state forward

2 c(t+1) = cost.fcn(s(t+1)); % calc cost distribution
```

The discounted cumulative cost is computed by iteration up to the horizon, propagating the state distribution forward and accumulating the discounted cost

$$f^{t} = f^{t-1} + \gamma^{t} c^{t}, \quad t = 1, ..., H, \text{ and } f^{t=0} = 0,$$
 (4)

where c^{t} is the instantaneous cost at time t and γ is the discount factor.

To accumulative costs and derivatives we again iterate up to the horizon, but now we need to keep track of the derivatives using the chain rule, as we move forward in time. When calling **propagated.m** (the derivative counterpart function of $\langle propagate.m 8 \rangle$) it computes the state distribution at time t from the state distribution at time t-1 and derivatives wrt both the state and the policy parameters:

$$s^{t} = g(s^{t-1}, p), \quad \frac{\partial s^{t}}{\partial s^{t-1}} \quad \text{and} \quad \frac{\partial s^{t}}{\partial p},$$
 (5)

where g is the *transition* function, which takes a distribution over states and a policy and returns the distribution over the next state. The derivatives are propagated forward using the chain rule

$$\frac{ds^{t}}{dp} = \frac{\partial s^{t}}{\partial s^{t-1}} \frac{ds^{t-1}}{dp} + \frac{\partial s^{t}}{\partial p}, \quad \text{where} \quad \frac{ds^{t=0}}{dp} = \mathbf{0}, \tag{6}$$

which is iterated forward in time in line 11 below. Finally the derivative of the cumulative discounted cost is the discounted cumulative derivatives in line 12 below

$$\frac{df^{t}}{dp} = \frac{df^{t-1}}{dp} + \gamma^{t} \frac{dc_{t}}{ds^{t}} \frac{ds^{t}}{dp} \quad \text{where} \quad \frac{df^{t=0}}{dp} = \mathbf{0}$$
 (7)

where f^t is the cumulative discounted cost up to time t, c^t is the instantaneous cost at time t and γ is the discount factor.

```
\langle compute\ cumulative\ costs\ and\ derivatives\ 7c \rangle \equiv
7c
                                                                             (5b)
      1 is = ctrl.is;
      2 ic = unwrap([is.m(1:D),is.s(1:D,1:D)]); % cost indices, depend on real vars
      3 sdp = cell(H+1); sdp{1} = zeros(ctrl.ns,ctrl.np);
                                                                           % init derivatives
      4 \text{ qdp} = \text{sdp}\{1\}(\text{is.s,:});
      5 dp = zeros(2,ctrl.np);
                                                 % first row = mean, second row = variance
      6 \text{ for } t = 1:H
                                                                      % iterate up to horizon
           [s(t+1), C, a(t), dsds, dsdp, dCds, dCdp] = propagated(s(t), dyn, ctrl);
           [c(t+1), dcds] = cost.fcn(s(t+1));
                                                            % cost and derivative wrt state
          cc.m = cc.m + gamma^t*c(t+1).m;
```

```
10
    cc.s = cc.s + gamma^{(2*t)*c(t+1).s};
11
    sdp\{t+1\} = dsds*sdp\{t\} + dsdp;
                                                                 % chain rule
12
    13
    % cross-covariance terms:
    dCdp = transposed(dCdp,C) + transposed(dCds,C)*sdp{t};
14
    qdp = prodd([],dCdp,[q,s(t+1).s]) + prodd(C',[qdp;sdp{t+1}(is.s,:)]);
15
16
    q = C'*[q, s(t+1).s];
17
    for j = 1:t % cross terms
      J = (j-1)*F+(1:D); % j'th column-block of q (real vars only)
18
19
      [cov, dcovdsj, dcovdst, dcovdq] = cost.cov(s(j), s(t+1), q(1:D,J)');
20
      cc.s = cc.s + 2*gamma^{(j+t-1)}*cov;
21
      dcov = dcovdsj*sdp{j}(ic,:) + dcovdst*sdp{t+1}(ic,:) + ...
22
        dcovdq*transposed(qdp(sub2ind2(F,1:D,J),:),D);
23
      dp(2,:) = dp(2,:) + 2*gamma^{(j+t-1)}*dcov;
24
    end
25 end
26 \text{ dcc.m} = dp(1,:);
27 \text{ dcc.s} = dp(2,:);
```

Two final subtleties consists in firstly, the derivatives of structures wrt to structures are represented as matrices (ie the structures have been vectorised), such that the products in line 11 and 12 are valid. This requires the structure dcds in line 12 to be vectorised (noting that c is a struct of two scalars). And secondly, note that in eq. (6) the state may contain fields in addition to s.m and s.s, but in eq. (7) only the fields s.m and s.s are considered (as the cost cannot depend directly on additional fields), and this is achieved by addressing directly those relevant indices using the predefined ctrl.is index structure in line 2.

2.3 Propagate

Propagate predicts the distribution over states at time t+1 given the distribution over states at time t. Predictions are made using a controller object to compute control actions, and then concatenating state and action information for input into a dynamics model.

```
\langle propagate.m \ 8 \rangle \equiv
1 function [s, C, a] = propagate(s, dyn, ctrl)
3 % Propagate the state distribution one time step forward.
5 % [s, C, a] = propagate(s, dyn, ctrl)
6 %
7 % s
                       state structure
8 %
              F x 1
                      mean vector
       m
9 %
              F x F
                       covariance matrix
       S
       ?
                       possibly other fields representing additional information
10 %
11 % dyn
                       dynamics model object
12 %
       D
                       dimension of the physical state
13 %
       Ε
                       dimension of predictions from dyn model
                      dynamics model function
14 %
              a
       pred
15 %
       pn
              E x 1
                      log std dev process noise
                      controller object
16 % ctrl
17 %
                      controller function
       fcn
18 %
       IJ
                      dimension of control actions
                      inverse input covariance times input-output covariance
19 % C
              F x F
20 % a
                       action structure
21 % m
              U x 1
                      mean vector
22 %
              U \times U
                      covariance matrix
23 %
```

```
24 % Copyright (C) 2008-2015 Carl Edward Rasmussen and Rowan McAllister 2015-06-05
26 F = length(s.m); D = ctrl.D; E = dyn.E; U = ctrl.U;
                                                                  % short hand names
27 Dz = F-D;
                                           % length of predicted information states
28 i = 1:D;
                                        % indices of physical state input variables
29 i = D+1:F;
                                                      % indices of information state
30 k = F + (1:U);
                                                        % indices of control actions
31 1 = \max(k) + (1:Dz);
                                           % indices of predicted information state
32 m = max([k,1]) + (1:E);
                                                       % indices of predicted states
33 ij = [i \ j]; ik = [i \ k]; kl = [k \ l]; ijkl = [ij \ kl];
                                                                         % short hand
34 \text{ o} = [ik(end-D+E+1:end) m 1];
                                                         % ind. to select next state
35 M = zeros(max(m),1); M(ij) = s.m; S = zeros(max(m)); S(ij,ij) = s.s;
37 [M(kl), S(kl,kl), A, s] = ctrl.fcn(s, dyn);
                                                     % control signal and inf state
38 q = S(ij,ij)*A; S(ij,kl) = q; S(kl,ij) = q';
                                                                % action covariances
40 [M(m), S(m,m), B] = dyn.pred(M(ik), S(ik,ik));
                                                      % compute distr of next state
41 S(m,m) = S(m,m) + diag(exp(2*dyn.pn));
                                                                 % add process noise
42 q = S(ijkl,ik)*B; S(ijkl,m) = q; S(m,ijkl) = q';
                                                            % next state covariances
44 C = [eye(F) A [eye(F,D) A(:,1:U)]*B];
                                                           % inv input var times cov
45 \% C_{exact_ZC} = [eye(F,D),A(:,U+1:end)]*C_gph;
47 \text{ s.m} = M(0); \text{ s.s} = (S(0,0)+S(0,0)')/2; C = C(ij,0);
                                                                 % select next state
48 a.m = M(k); a.s = (S(k,k)+S(k,k)')/2;
                                                                    % control action
```

On line 37 the controller outputs the matrix A. A which is the covariance between the input variable (the state, distributed as a Gaussian of mean s.m and variance s.s) and the output variable (a concatenation of the action variable and the state filter prediction variables, distributed as a Gaussian of mean M(kl) and variance S(kl,kl)) pre-multiplied by the inverse variance of the input variable. I.e.

$$A = V[s^{t}]^{-1}C[s^{t}, \{u^{t}, s_{F-D+1:F}^{t+1}\}],$$
(8)

where s^t is the state at time t, u^t is the action taken at time t, and $s^{t+1}_{F-D+1:F}$ is the predicted filter state variables (whose indexes are are [F-D+1,F]). On line 40 the dynamics model outputs the matrix B. Similarly, B is an input-output covariance matrix pre-multiplied by the inverse of input variance. We have

$$B = V[\{s^t, u^t\}]^{-1}C[\{s^t, u^t\}, s^{t+1}], \tag{9}$$

where s^t is the state at time t and u^t is the action taken at time t resulting in state s^{t+1} at time t+1. The use of an pre-multiplied variance inverse of the input variable has the nice property that the covariance between one variable and any ancestor is simply computed with the product of the ancestor's variance and all such C terms along the path in between.

2.4 Rollout

We define a sampled trajectory of a system's possible evolution up to horizon H as a *rollout*. The system transitions from (point mass) state s^t to (point mass) state s^{t+1} as the system observes point observations y^t and applies point control signals u^t at each point in time t. If the state struct s does not have a variance field (i.e if s.s does not exist), then subroutines will assume they are called by **rollout** opposed to **propagate**.

```
9 ⟨rollout.m 9⟩≡
1 function [data, latent, L] = rollout(start, ctrl, H, plant, cost, verb)
2 % Compute a state trajectory using an ode solver (and any additional dynamics)
3 % from a particular starting state with either a particular policy or random
4 % actions.
```

```
6 % [data, latent, L] = rollout(start, ctrl, H, plant, cost, verb)
 7 %
 8 % start nX x 1 vector containing start state (without controls)
9 % ctrl controller structure
10 % fcn @ function implementing the controller
11 % init @ function initialising controller's filtered state
12 % policy structure
11 % init @ policy policy structure
13 % fcn @ policy function
14 % p p parameter structure (if empty use random actions)

1 vector of control input saturation values
17 % plant
                           the dynamical system structure
19 % augment (opt) augment state using a known mapping
20 % constraint (opt) stop rollout if violated
21 % dyno indices for states passed to cost
22 % noise observation noise
23 % odei indices for states passed to the ode solver
24 % poli indices for states passed to the policy
25 % cost cost object
26 % yerb
18 % augi
                          (opt) indices for states passed to augment function
                          verbosity level
26 % verb
27 %
28 % data
                 data struct
28 \% data data struct 29 \% state H+1xnX state matrix
30 \% action H x nU action matrix
31~\% L loss incurred at each timestep (1 by H)
32 % latent matrix of latent states (H+1 by nX)
34 % Copyright (C) 2012-2015 Carl Edward Rasmussen and Rowan McAllister 2015-06-05
35
36 clear odestep;
                                               % clear persistent old action function
38 else plant.augment = @(x)[]; augi = []; end
39 calc_loss = nargout > 2;
40 if nargin < 6; verb = 0; end
41 D = ctrl.D; E = ctrl.E; F = ctrl.F; U = ctrl.U;
42 N = length(start); odei = plant.odei;
43 latent = zeros(H+1, N); y = NaN(H+1, N); L = zeros(1, H+1); u = zeros(H, U);
44 obs_noise = @()(randn(1,E)*chol(plant.noise));
45
46 latent(1,:) = start;
                                                                            % initialise
47 y(1,1:D-E) = latent(1,1:D-E);
48 y(1,D-E+1:D) = latent(1,D-E+1:D) + obs_noise(); % add noise to observations
49 if calc_loss; L(1) = cost.fcn(struct('m',latent(1,1:D)')).m; end
51 s.m = y(1,1:D)'; s = ctrl.reset_filter(s); % reset filter if exists
53 for i = 1:H % ------ run ROLLOUT
54 % Test constraints and stop rollout if violated ------
    if isfield(plant, 'constraint') && plant.constraint(latent(i,:))
56
     H = i-1; if verb; disp('state constraints violated...'); end; break;
57
    end
58
59
    % Apply policy ------
    s.m(1:D) = y(i,1:D)'; % receive an observation
60
61 [uzm, \sim, \sim, s] = ctrl.fcn(s);
62
     u(i,:) = uzm(1:U); % action 'u' component of uzm
```

```
s.m(D+1:F) = uzm(U+1:end); % predicted filter 'zm' component of uzm
63
64
65
    latent(i+1,1:D-E) = [latent(i,U+E+1:D), u(i,:)];
66
    latent(i+1,odei) = odestep(latent(i,odei), u(i,:), plant);
67
    y(i+1,1:D-E) = [y(i,U+E+1:D), u(i,:)];
69
    y(i+1,D-E+1:D) = latent(i+1,D-E+1:D) +obs_noise(); % TODO: add process noise?
70
    % Compute Cost ------
71
72
    if calc_loss; L(i+1) = cost.fcn(struct('m',latent(i+1,1:D)')).m; end
74 if verb; disp(['Trial lasted ',num2str(floor(H)),' steps']); end
76 data.state = y(1:H+1,:); data.action = u(1:H,:);
77 latent = latent(1:H+1,:); L = L(1,1:H+1);
                                                   % trim any trailing zeros
78
79 function xa = augment(x, plant)
80 xa(plant.odei) = x;
81 xa(plant.augi) = plant.augment(xa);
```

3 Controllers

Controller classes act as wrappers to policy functions. The superclass Ctrl.m defines the common functions a controller object supports, discussed Sec. 3.1. Then follows two sections on child classes that inherit Ctrl.m; the 'No Filter' controller CtrlNF.m in Sec. 3.2 and the 'Bayes Filter' controller CtrlBF.m Sec. 3.3.

3.1 Ctrl

```
11
     \langle Ctrl.m \ 11 \rangle \equiv
     1 classdef Ctrl < handle
         %CTRL, controller superclass, which ctrlBF and ctrlNF inherit.
     3
     4
         % Ctrl Properties:
                                     (optional) call this function with the action
     5
         %
             actuate - @
     6
         %
             angi
                                     indicies for vars treated as angles (sin/cos rep)
     7
         %
            D
                                     number of real-world state variables
     8
         % dyn
                                     dynamics model object (only used by CTRLBF class)
     9
         % E
                                     number of physical state variables
    10
             F
                                     number of state variables (real + info vars)
         %
         %
                                     state index structure
    11
            is
    12
         %
                                     number of policy parameters
             np
                                     state parameters (including filter)
    13
         %
             ns
                           - D x D non-log observation noise
    14
         %
             on
    15
         %
             onp

    D x D non-log obs. noise (only last E vars non zeros)

    16
         %
             poli
                                     indicies for the inputs to the policy
             policy
    17
         %
                                     policy struct
    18
         %
             U
                                     number of control outputs
    19
         %
         % Ctrl Methods:
    20
    21
         % build_state_index - (private) builds 'is' field to index state-struct
    22
         % clear_filter - clears all filter variables (if they exist)
    23
        % Ctrl

    constructor

    24
        % fcn
                               - main function, computes control signal
         % random_action - outputs a random control, independent of state
% reset_filter - resets state fields {zs,zc,v} to a broad prior
    25
    26
```

```
27
                           - sets the dynnamics model object
        set_dynmodel
28
                               sets the N-Markov observations noise
    %
        set_on
                          - sets the policy optimisation settings
29
    %
        set_policy_opt
                               sets the policy parameters
30
        set_policy_p
                          -
31
    %
32
    % See also CTRLBF.M, CTRLNF.M.
    % Copyright (C) 2015 by Carl Edward Rasmussen and Rowan McAllister 2015-06-05
33
34
35
    properties (SetAccess = private)
36
      actuate
37
      angi
38
39
      dyn
40
      Ε
41
      F
42
      is
43
      np
44
      ns
45
      on
46
      onp
47
      poli
48
      policy
49
      U
50
    end
51
52
    methods
53
54
      % Constructor
55
      function self = Ctrl(D, E, policy, angi, poli, actuate)
56
        % CTRL is the super-class controller constructor
57
        % ctrl = Ctrl(D, E, policy, angi, poli, actuate)
58
59
60
        % INPUTS:
                              number of state variables
61
        % D
62
        %
                              (or a ctrl object to be copied)
63
        % E
                              number of predicted state variables
64
        % policy
                              policy struct
65
                              policy function
        % fcn
                       a
                              maximum control output magnitudes
66
        %
              maxU
                              optimisation structure
67
        %
            opt
68
        %
              fh
                              figure handle for minimize() to display to
69
               length
                              how many optimisations steps
        %
70
        %
                method
        %
                verbosity
71
72
                              indicies for variables treated as angles
        % angi
                              indicies for the inputs to the policy
73
            poli
        %
74
                              function to actuate calculated action
            actuate @
75
76
        % Special case input:
        % Another controller object might be the first (and only) input.
77
        % This allows easy translations from one controller type to the next
78
79
        if ~isnumeric(D); assert(nargin < 2); ctrl = D;</pre>
          D = ctrl.D;
80
81
          E = ctrl.E;
82
          policy = ctrl.policy;
83
          angi = ctrl.angi;
84
          poli = ctrl.poli;
```

```
85
            actuate = ctrl.actuate;
 86
            self.on = ctrl.on;
 87
            self.onp = ctrl.onp;
 88
            self.dyn = ctrl.dyn;
 89
          end
 90
 91
          self.D = D;
          self.E = E;
 92
 93
          self.policy = policy;
 94
          if ~isfield(policy,'opt'); self.policy.opt = ...
 95
              struct('length',-1000,'method','BFGS','MFEPLS',20,'verbosity',3); end
          if ~isfield(self.policy.opt,'fh'); self.policy.opt.fh = 1; end
 96
 97
          self.U = length(policy.maxU);
 98
          self.build_state_index();
99
          if ~isfield(policy,'type'); policy.type = ''; end
100
          if isfield(self.policy,'p');
101
            self.np = length(unwrap(self.policy.p));
102
          end
103
104
          if exist('angi','var'); self.angi = angi; else self.angi = []; end
105
          if exist('poli','var'); self.poli = poli; else self.poli = 1:self.D; end
106
          if exist('actuate','var') && ~isempty(actuate); self.actuate=actuate; end
107
108
109
        function [uM,uS,uC,s] = fcn(self, s)
110
          % CTRL.FCN is the main function to output control. Sub-classes will
          % override this function.
111
112
113
          % [uM, uS, uC, s, duMds, duSds, duCds, dsds, duMdp, duSdp, duCdp, ...
114
              dsdp] = CTRL.FCN(s,propdyn)
115
         %
116
         % self
                                   controller structure
117
         %
                                   function to actuate calculated action
             actuate
118
          %
             angi
                                   indices of angular variabels
119
                                   controller's dynamics model (for CtrlBF only)
         %
             dyn
            E
120
         %
                                   number of predictive state variables
121
                        D \times D
                                   observation noise
         %
            on
122
         % onp
                        D \times D
                                   observation noise (only last E vars non-zero)
123
         %
                                   indices of policy input
            poli
124
         %
              policy
                                   policy structure
125
         %
                fcn
                         a
                                   policy function
126
                                   number of control outputs
         %
127
         % s
                                   state structure
128
                         F x 1
          %
            m
                                   state mean
129
          %
                         F x F
                                   state variance
            S
130
          %
             V
                    (F-D) x (F-D) filter variance
131
                                   propagates's dynmodel (for CtrlBF)
         % propdyn
                        a
132
         % M
                    (U+D) \times 1
                                   control signal mean vector
133
         % S
                    (U+D) x (U+D) control signal variance matrix
134
         % C
                         F x (U+D) input-output covariance matrix
         % dMds
135
                    (U+D) \times S
                                   derivatives of outputs wrt input state struct
136
         % dSds (U+D)^2 \times S
137
         % dCds F*(U+D) \times S
         % dsds
                         S \times S
                                   ouput state derivative wrt input state
138
139
         % dMdp
                    (U+D) \times P
                                   P is the total number of parameters is the policy
140
         % dSdp (U+D)^2 \times P
141
          % dCdp F*(U+D) \times P
142
                         S x P
         % dsdp
                                   ouput state derivative wrt policy parameters
```

```
143
144
         % See also CTRLBF.FCN, CTRLNF.FCN.
145
         if strcmp(self.policy.type, 'random')
146
           uM = self.random_action();
147
            uS = zeros(self.U); uC = zeros(self.D,self.U);
148
          end
       end
149
150
       % Filter Clearer
151
152
        function s = clear_filter(self, s)
         % Clears all filter variables (if they exist).
153
154
         % s = CTRL.CLEAR_FILTER(s)
         % s: state struct
155
         i = 1:self.D;
156
157
          s.m = s.m(i);
158
          if isfield(s,'s'); s.s = s.s(i,i); end
159
          if isfield(s,'v'); s = rmfield(s,'v'); end
160
       end
161
162
       function u = random_action(self)
163
         % u = CTRL.RANDOM_ACTION(), outputs a random action
164
         % u:
                  U x 1 control action
165
         u = self.policy.maxU.*(2*rand(1,self.U)-1);
166
       end
167
168
       function s = reset_filter(~, s)
169
         % s = CTRL.RESET_FILTER(s), does nothing unless overwritten
170
         % s: state struct
         % See also CTRLBF.RESET_FILTER
171
172
        end
173
174
       function set_dynmodel(self, dyn)
175
         % CTRL.SET_DYNMODEL(dyn), for updating controller's dynmodel
176
              dyn: dynamics model object
177
         % Note: dyn property only ever used by CTRLBF
178
          self.dyn = dyn;
179
        end
180
181
       function set_on(self, onE)
182
         % CTRL.SET_ON(onE), sets observation noise
183
             onE: E x 1, log observation noise (physical state-variables only)
184
          assert(self.E == length(onE));
         onD = [-inf(self.U,1); onE(:)];
185
186
          onD = repmat(onD,ceil(self.D/(self.U+self.E)),1);
187
          onD = onD(end-self.D+1:end);
188
          self.on = diag(exp(2*onD));
189
         p = self.D-self.E+1:self.D;
190
          self.onp = 0*self.on; self.onp(p,p) = self.on(p,p);
191
192
193
        function set_policy_p(self, p)
194
         % CTRL.SET_POLICY_P(p), updateds the policy parameters
195
              p: policy parameter struct
196
          self.policy.p = p;
197
          self.np = length(unwrap(p));
198
       end
199
200
       function set_policy_opt(self, opt)
```

```
% CTRL.SET_POLICY_OPT(p), updateds the policy optimisation settings
201
202
                           policy optimisation settings
                  opt:
203
            self.policy.opt = opt;
204
         end
205
206
       end
207
       methods (Access = private)
208
209
          \langle function\text{-}build\text{-}state\text{-}index \ 15 \rangle
210
       end
211
212 end
```

PILCO stores derivatives of any object as matrices, where rows represent vectorised dependent variables, and columns represent vectorised independent variables. To vectorise any object we use the unwrap.m function. The state distribution is stored as a structure. Thus when handling state derivatives, we need to keep track of which elements in the native structure correspond to which index in the vectorisation of that structure. We do so by using a *state index* structure, which has the exact same form of the state structure, except each element within the index structure is a unique integer, corresponding to that element's index when in vectorised form. The state index is built using \(\frac{\tau tunction-build-state-index}{\text{15}} \):

```
\langle function\text{-}build\text{-}state\text{-}index \ 15 \rangle \equiv
15
                                                                             (11)
      1 function build_state_index(self)
          % CTRL.BUILD_STATE_INDEX(), computes internal field 'is' - a state struct
      3
          % whose members are a states members' indexes. Requires subclass to
          % implement function CTRL.RESET_FILTER
      4
      5
          s.m = nan(self.D,1);
          s = self.reset_filter(s); % generates possible filter variables
      7
          s.s = nan(length(s.m));
          if isfield(s,'reset'); s = rmfield(s,'reset'); end
      8
          self.ns = length(unwrap(s));
      9
     10
          self.is = rewrap(s,1:self.ns);
          self.F = length(s.m);
     11
     12 end
```

3.2 CtrlNF

The most straightforward controller class is one which inputs (noisy) signals from sensors directly into the policy function. The 'No Filter' controller class **control/CtrlNF** implements exactly this.

The process of controlling a dynamical system using CtrlNF.m is depicted in Fig. 1. It is assumed the latent system state X_t is observed by imperfect sensors as Y_t , where the random variable Y_t is equal to X_t plus some Gaussian noise:

$$Y_t \sim X_t + \varepsilon_t,$$
 (10)

$$\epsilon_{t} \stackrel{\text{iid}}{\sim} \mathcal{N}(0, \Sigma_{\epsilon}).$$
(11)

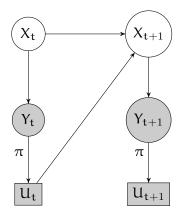


Figure 1: Directed graphical model of a system controller with No Filter from timestep t to timestep t+1. A dynamical system begins in state X_t , which sensors observe noisily as Y_t . The controller uses Y_t to decide control signal U_t via policy function π . Finally, the control signal U_t is applied to the system, resulting in new state X_{t+1} .

```
\langle CtrlNF.m \ 16a \rangle \equiv
16a
       1 classdef CtrlNF < Ctrl & handle</pre>
       3
           % Controller with No Filter. The state is given either as a point s.m or as a
       4
           % distribution N(s.m,s.s). First augment with trignometric functions if
           % necessary, then call the policy, and finally (optionally) call the actuate
           % function. There is no filter, so no updates to any state structure filter
       6
       7
           % fields are required.
       8
       9
           % See also CTRL.M, CTRLNFT.M.
      10
           % Copyright (C) 2015 by Carl Edward Rasmussen and Rowan McAllister 2015-06-02
      11
      12
           methods
      13
      14
              % Constructor
      15
              function self = CtrlNF(varargin)
      16
                % CTRLNF.CTRLNF is the sub-class constructor of CTRL.CTRL
      17
                % For help, see also Ctrl.Ctrl
      18
                self@Ctrl(varargin{:});
                                                                       % Super Ctrl constructor
      19
              end
      20
      21
              % Main function
      22
              function [uM, uS, uC, s, duMds, duSds, duCds, dsds, ...
      23
                  duMdp, duSdp, duCdp, dsdp] = fcn(self, s, ~)
      24
                % For help, see also CTRL.FCN
      25
      26
                (initialise 16b)
      27
                \langle augment \ state \ with \ trig \ variables \ 17 \rangle
      28
                ⟨compute control signal 18⟩
      29
              end
      30
      31
           end
      32
      33 end
16b
       \langle initialise \ 16b \rangle \equiv
                                                                             (16a)
       1 % CtrlNF only operates on a noisy version of the state:
```

2 D = self.D; DD = D*D;

```
3 s = self.clear_filter(s);
                                       % clear any filter variables in state
4 assert(length(s.m) == D);
5 if isfield(s,'s')
    sy = s.s + self.on;
                                  % propagate mode (distribution of states)
7 \; {\it else}
    sy = zeros(D);
                                    % rollout mode (point-mass state sample)
8
9 end
11 angi = self.angi; poli = self.poli; A = length(angi);
12 derivativesRequested = nargout > 4;
13 ns = self.ns; is = self.is;
14 D1 = D + 2*A;
15 i=1:D;
16 M = zeros(D1,1); M(i) = s.m; S = zeros(D1); S(i,i) = sy;
17 if derivativesRequested
    idx = @(i,j,I) bsxfun(@plus, I*(i'-1), j);
19
    Mds = zeros(D1,ns); Mds(i,is.m) = eye(D);
    Sds = zeros(D1*D1,ns); Sds(:,is.s) = kron(Mds(:,is.m),Mds(:,is.m));
20
21
    dsds = zeros(self.ns); dsdp = zeros(ns,self.np);
    dsds(is.m,is.m) = eye(D); dsds(is.s(:),is.s(:)) = eye(D*D);
23
    dsds(is.s,is.s) = symmetrised(dsds(is.s,is.s),[1,2]);
24 end
```

The state distribution s, is then extended the sine and cosine each angle variable (indexed by angi) using gTrig.m. The gTrig.m function also outputs covariance information C_{gtrig} , which relates to the input-output covariance except with pre-multiplicative input variance:

$$\mathbb{C}[Y_t, Y_t^{\circ}] = \mathbb{V}[Y_t] C_{\mathfrak{atrig}}[Y_t, Y_t^{\circ}] \tag{12}$$

```
17
      \langle augment \ state \ with \ trig \ variables \ 17 \rangle \equiv
                                                                         (16a)
      1 % augment state with trig functions
      2 i = 1:D; k = D+1:D1;
      3 if ~derivativesRequested
          [M(k), S(k,k), cg] = gTrig(M(i), S(i,i), angi);
      5 else
      6
          kk = idx(k,k,D1); ik = idx(i,k,D1); ki = idx(k,i,D1);
          [M(k), S(k,k), cg, Mds(k,is.m), Sds(kk,is.m), cgdm, ...
            Mds(k,is.s), Sds(kk,is.s), cgds] = gTrig(M(i), S(i,i), angi);
      8
      9
          qdm = prodd(S(i,i),cgdm);
          Sds(ik,is.m) = qdm; Sds(ki',is.m) = qdm;
     10
     11
          qds = prodd(S(i,i),cgds) + prodd([],'eye',cg);
     12
          Sds(ik,is.s) = qds; Sds(ki',is.s) = qds;
     14 q = S(i,i)*cg; S(i,k) = q; S(k,i) = q';
```

The policy, which may be a function of both the regular state variables, and augmented state variables is then computed:

$$U_t \leftarrow \pi(\{Y_t, Y_t^{\circ}\}) \tag{13}$$

The non-trivial part is computing output uC which is covariance between state input and control output $\mathbb{C}[X_t, U_t]$. We begin noting:

$$\mathbb{C}[\{Y_t, Y_t^{\circ}\}, U_t] = \mathbb{V}[\{Y_t, Y_t^{\circ}\}] C_{poli}[\{Y_t, Y_t^{\circ}\}, U_t]$$

$$\tag{14}$$

$$\begin{bmatrix} \mathbb{C}[Y_t, U_t] \\ \mathbb{C}[Y_t^{\circ}, U_t] \end{bmatrix} = \begin{bmatrix} \mathbb{V}[Y_t] & \mathbb{C}[Y_t, Y_t^{\circ}] \\ \mathbb{C}[Y_t^{\circ}, Y_t] & \mathbb{V}[Y_t^{\circ}] \end{bmatrix} C_{poli}[\{Y_t, Y_t^{\circ}\}, U_t]$$
(15)

$$\therefore \mathbb{C}[Y_t, U_t] = \left[\mathbb{V}[Y_t], \mathbb{C}[Y_t, Y_t^{\circ}] \right] C_{poli}[\{Y_t, Y_t^{\circ}\}, U_t]$$
 (16)

$$= \mathbb{V}[Y_t][I, C_{gtrig}[Y_t, Y_t^{\circ}]]C_{poli}[\{Y_t, Y_t^{\circ}\}, U_t]$$

$$\tag{17}$$

Thus,

18

$$\mathtt{uC} = \mathbb{C}[X_t, U_t] = \mathbb{C}[X_t, Y_t] \mathbb{V}[Y_t]^{-1} \mathbb{C}[Y_t, U_t] \tag{18}$$

$$= \mathbb{V}[X_t][I, C_{\mathfrak{gtrig}}[Y_t, Y_t^{\circ}]]C_{\mathfrak{poli}}[\{Y_t, Y_t^{\circ}\}, U_t]$$
 (19)

```
\langle compute\ control\ signal\ 18 \rangle \equiv
                                                                     (16a)
1 % compute control signal
2 if ~derivativesRequested
     [uM, uS, uC] = self.policy.fcn(self.policy, M(poli), S(poli,poli));
4 else
     [uM, uS, uC, mdm, sdm, cdm, mds, sds, cds, duMdp, duSdp, duCdp] = ...
6
       self.policy.fcn(self.policy, M(poli), S(poli,poli));
8 if isfield(self, 'actuate'), self.actuate(uM); end  % actuate controller
10 \text{ ec} = [\text{eye}(D) \text{ cg}]; \text{ ecp} = \text{ec}(:,\text{poli});
11 if derivativesRequested
     poli2 = idx(poli,poli,D1); ii = sub2ind2(D1,i,i);
     duMds = mdm*Mds(poli,:) + mds*Sds(poli2,:);
13
14
     duSds = sdm*Mds(poli,:) + sds*Sds(poli2,:);
15
     duC = cdm*Mds(poli,:) + cds*Sds(poli2,:);
16
17
     decp = [zeros(DD,ns); cgdm*Mds(i,:) + cgds*Sds(ii,:)];
     decp = decp(sub2ind2(D,1:D,poli),:);
18
19
     duCds = prodd(ecp,duC) + prodd([],decp,uC);
20
21
     duCdp = prodd(ecp,duCdp);
22
23
     duCds(:,is.s) = symmetrised(duCds(:,is.s),2);
24
     duMds(:,is.s) = symmetrised(duMds(:,is.s),2);
25
     duSds(:,is.s) = symmetrised(duSds(:,is.s),2);
26 end
27 \text{ uC} = ecp*uC;
```

3.3 CtrlBF

CtrlBF.m implements control using Bayesian filtering. The *state* of the control system using filtering contains two parts:

1. the latent state x,

2. the filter state distribution $b \sim \mathcal{N}(z, V)$.

Thus, the *state distribution* is in principle a distribution over the random variables, x, z and V. However, as an approximation we are going to assume that the distribution on the variance V is just a delta function (ie, that the variance is some fixed value). Assuming further that the state distribution is Gaussian

$$\begin{bmatrix} x \\ z \end{bmatrix} \sim \mathcal{N} \left(\begin{bmatrix} m_x \\ m_z \end{bmatrix}, \begin{bmatrix} \Sigma_x & \Sigma_{xz} \\ \Sigma_{zx} & \Sigma_z \end{bmatrix} \right). \tag{20}$$

A Bayes filter (BF) maintains a belief-posterior on x, denoted B_{tt} , conditioned on all information available thus far: the entire history of the system observations $y_{1:t}$ and applied control signals $\mathfrak{u}_{1:t-1}$. Conditioning on more information than the current observation y_t yields a more informed (and smooth) estimate of x_t . Being a function of all observations, B_{tt} is less susceptible to the noise injected into the most recent observation y_t , and consequently the controller's input is much smoother. To maintain B_{t+1} the BF makes two recursive updates per timestep:

- 1. Update step: Compute B_{tt} using prior belief $B_t = p(x_t)$ and observation likelihood $\mathcal{L}(x_t|y_t) = p(y_t|x_t)$,
- 2. Predict step: Compute B_{t+1} by mapping updated belief B_{tt} through transition model $p(x_{t+1}|b_{tt},u_t)$.

A directed graphical model of a Bayes filter is shown Fig. 3.3:

