

AA203 Optimal and Learning-based Control

Lecture 18

Model-based Policy Learning

Autonomous Systems Laboratory
Daniele Gammelli

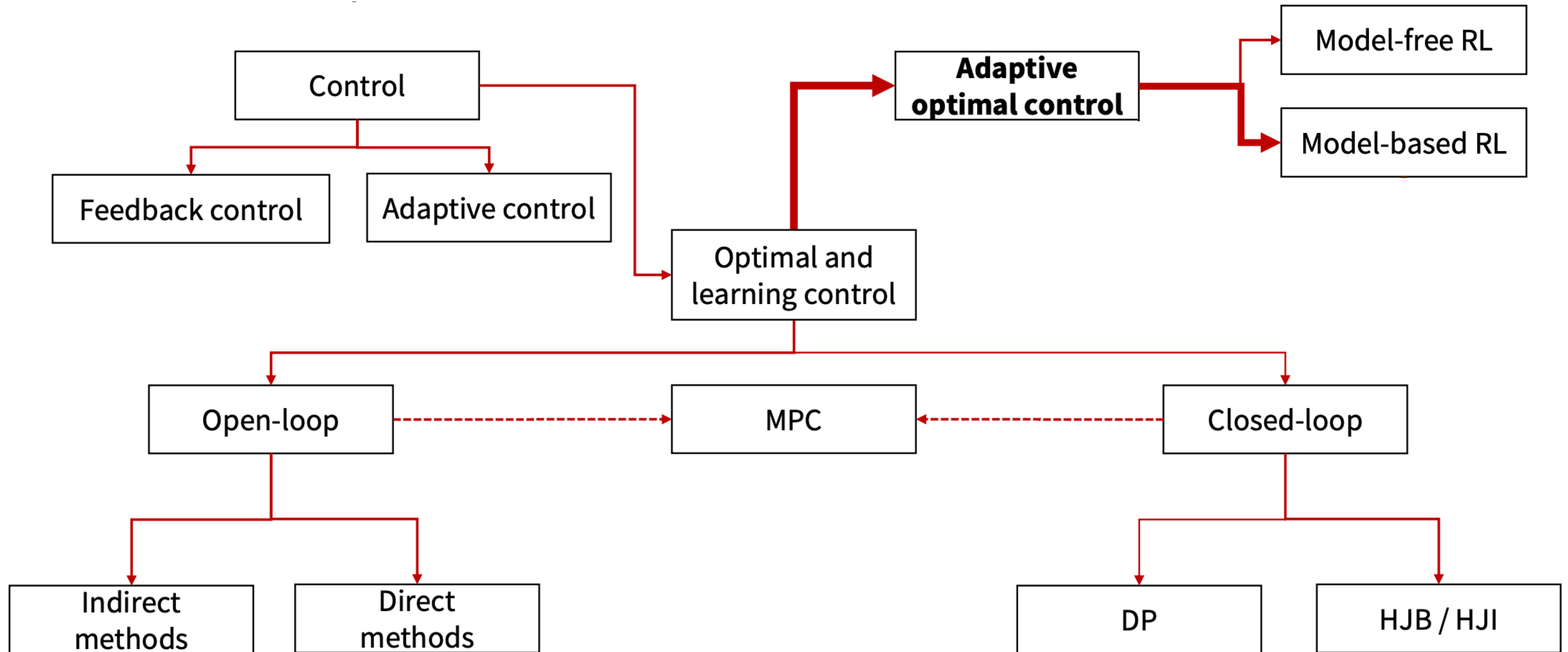


Stanford University



**Autonomous Systems Laboratory
Stanford Aeronautics & Astronautics**

Roadmap



Recap: Model-based RL

- In model-free RL, we discussed different approaches to solve unknown MDPs directly from experience via **policy / value-function learning**
- In model-based RL, we aim to (1) **estimate an approximate model** of the dynamics, and (2) **use it for control**

Approach 1: “learn a model $p(x_{t+1} | x_t, u_t)$ from experience and use it to plan”

1. Run base policy $\pi_0(u_t | x_t)$ in the environment (e.g., random policy, exploration policy) and collect dataset of transitions $\mathcal{D} = \{(x_t, u_t, x_{t+1})_i\}$

2. Fit dynamics model to data to minimize error (or equivalently, maximize (log) likelihood)

$$\theta^* = \arg \min_{\theta} \sum_i \left\| f_{\theta}(x_t, u_t) - x_{t+1} \right\|^2$$

YES

NO

Sys. ID

Distribution mismatch

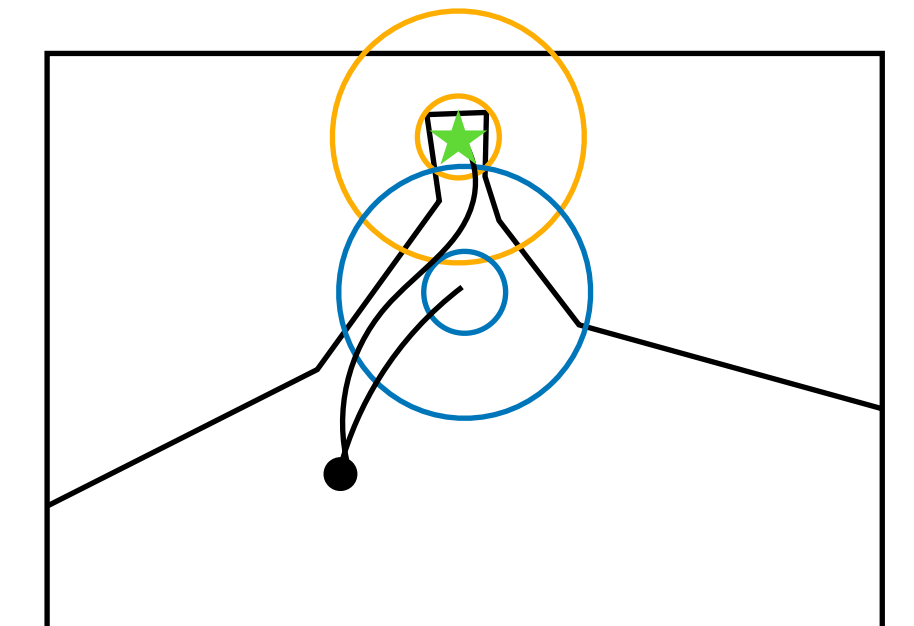
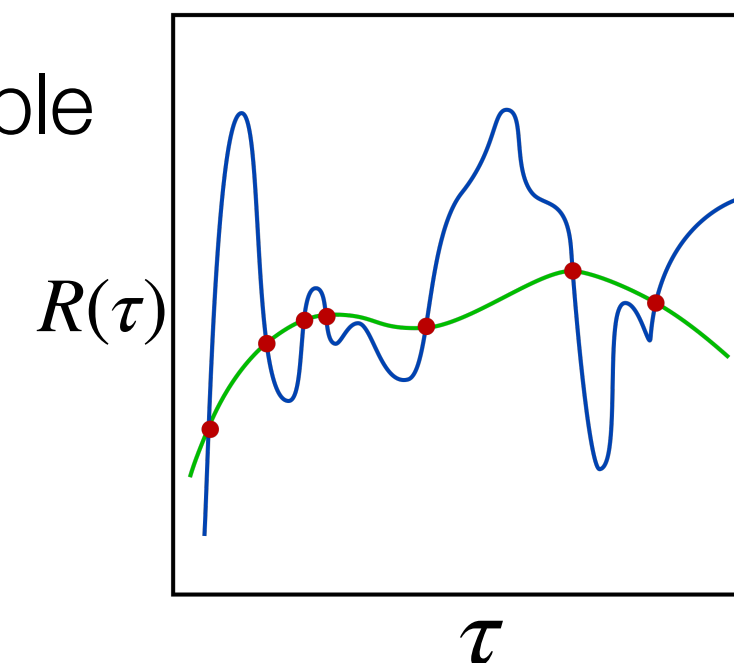
3. Use the learned model to plan a sequence of actions

Exploitation of errors

Problem: we'll likely *erroneously* exploit our model where it is less knowledgeable

(Possible) Solution: consider how “certain” we are our about the prediction

This allows us to reason in terms of expectations under our model

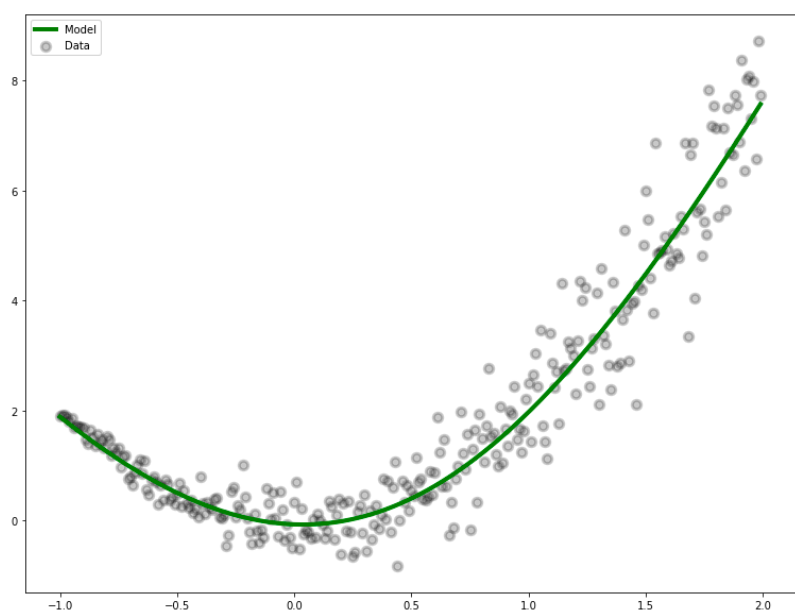


Recap: Model-based RL

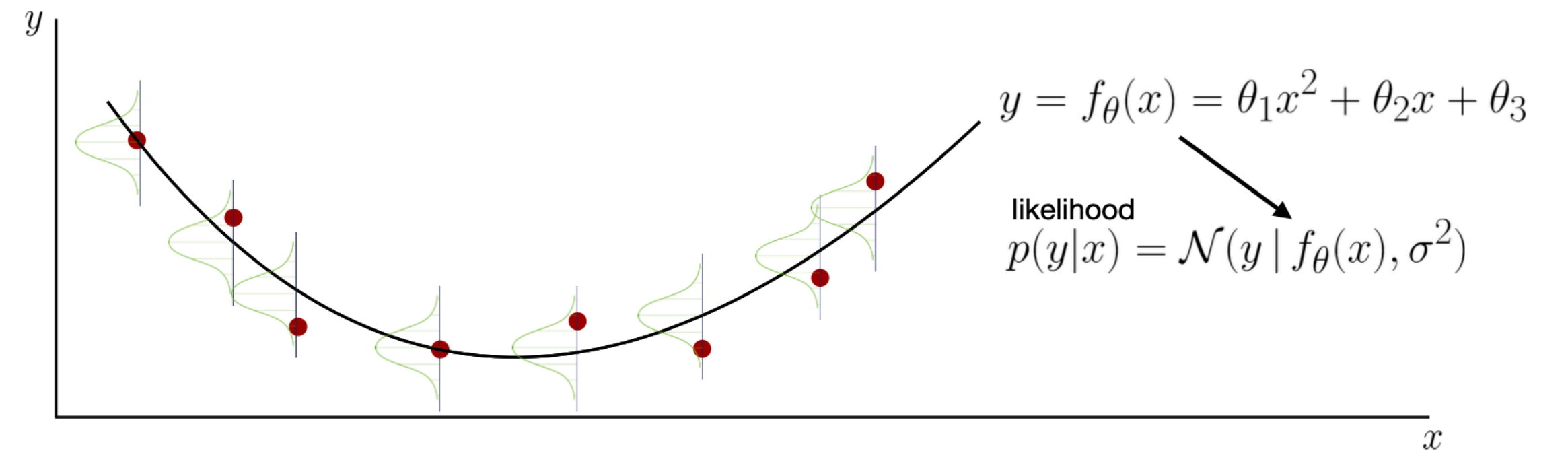
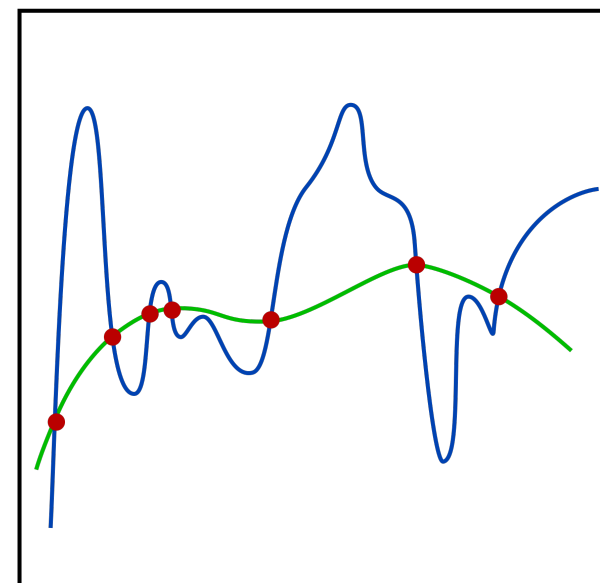
Uncertainty estimation

- Learning from a probabilistic standpoint (i.e., deterministic vs probabilistic predictions)
- The importance of estimating *model/epistemic uncertainty*

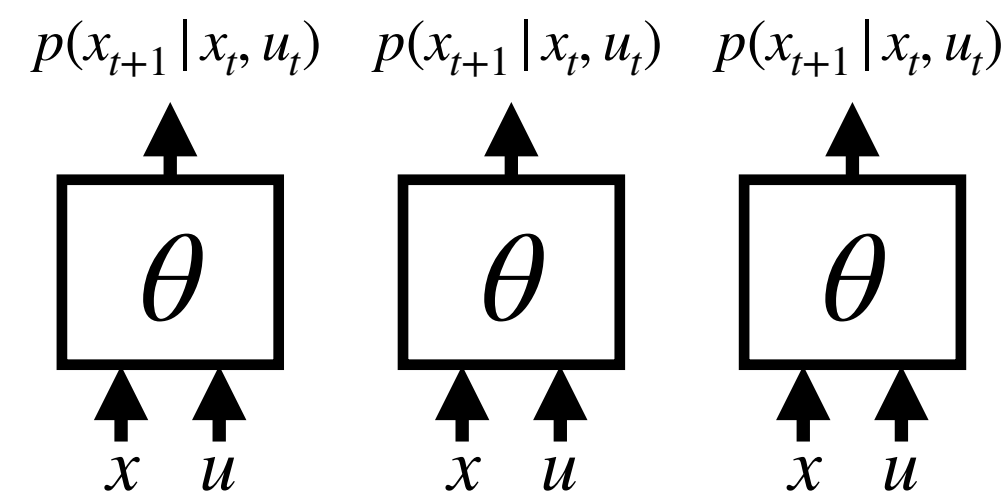
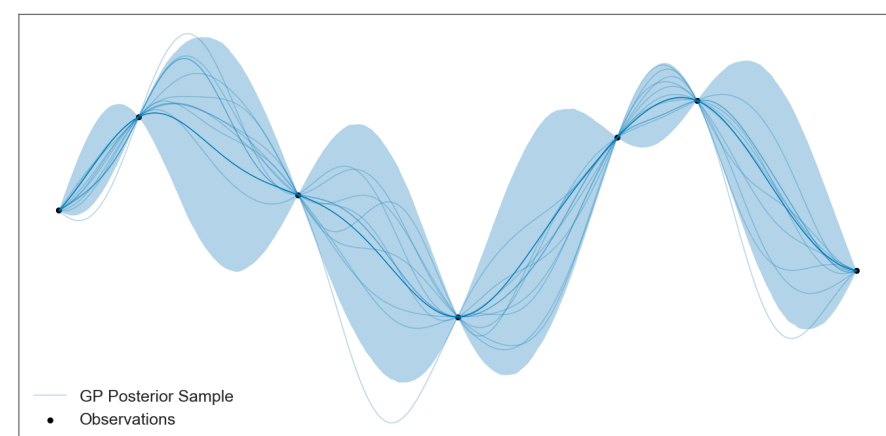
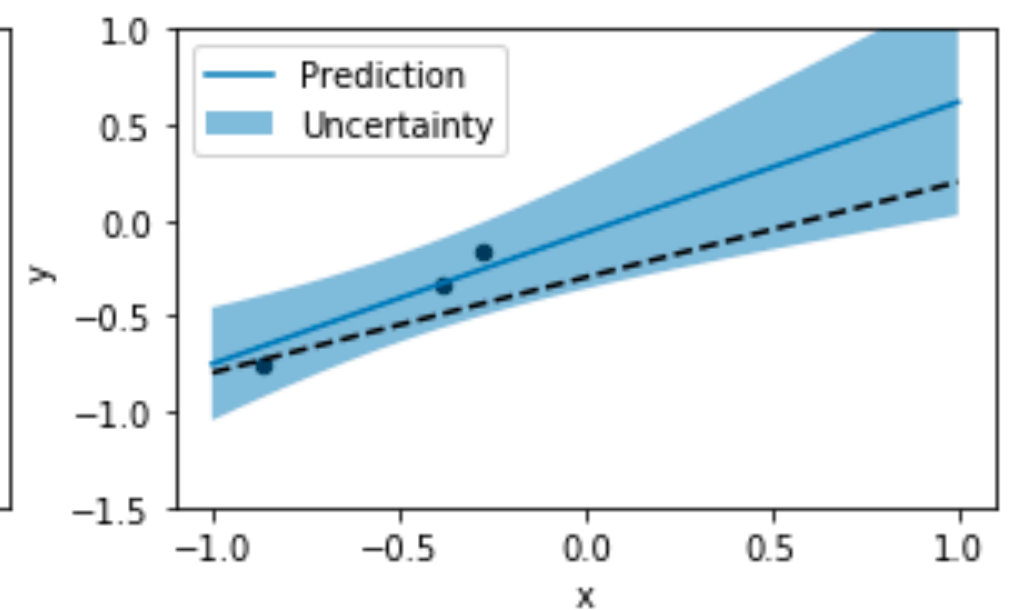
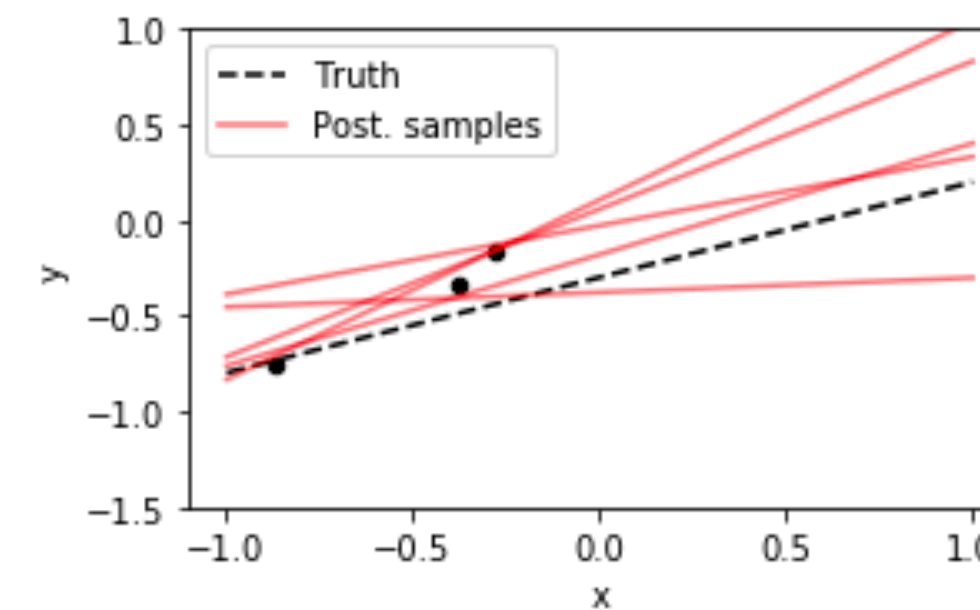
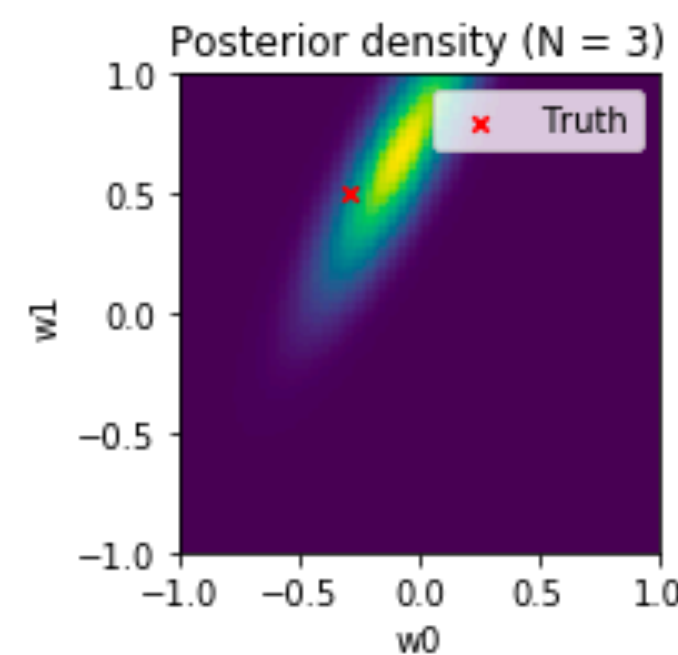
Aleatoric uncertainty



Epistemic uncertainty



- A structured way to represent uncertainty over a parametric model is through a **posterior distribution** over the parameters $p(\theta | \mathcal{D})$

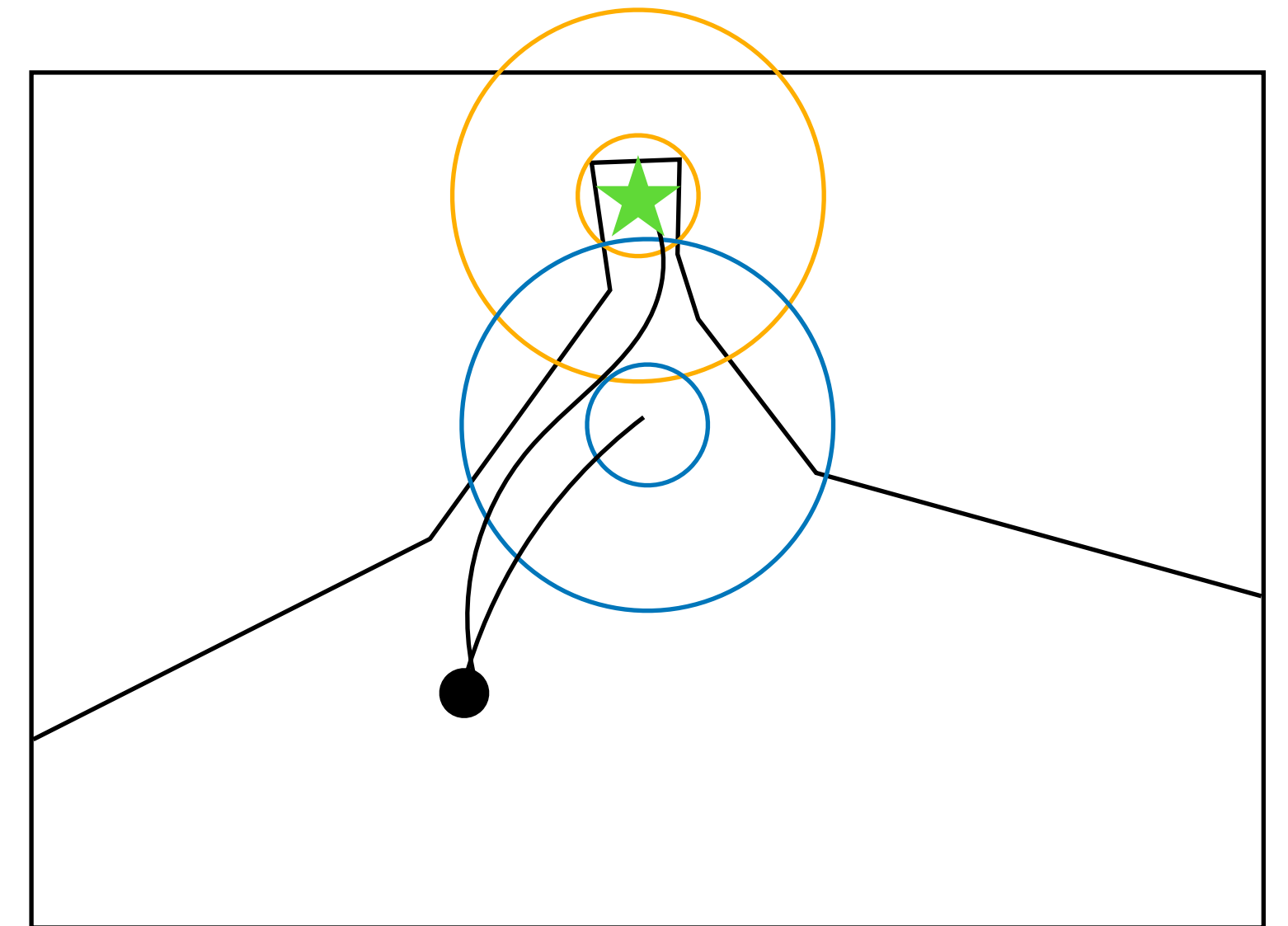


- Two examples:
 - Gaussian Processes (accurate; expensive; limited expressivity)
 - Ensembles (approximate; simple; high-capacity NNs)

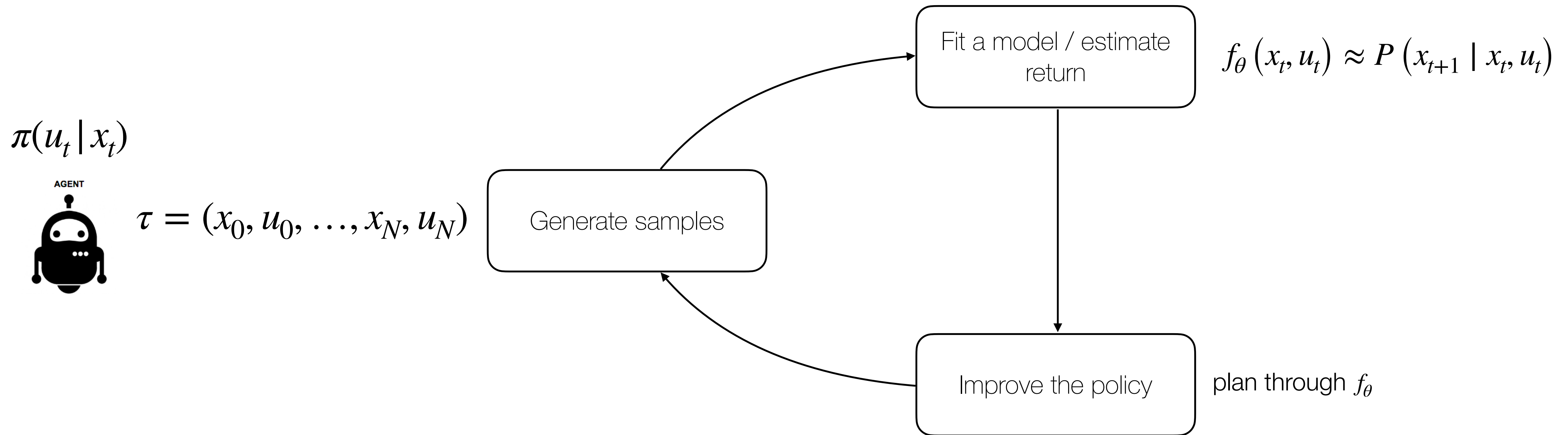
Recap: Model-based RL

- How do we use this in planning? A possible idea is the following:
- Given a candidate action sequence u_1, \dots, u_T :
 1. Sample $\theta_i \sim p(\theta | \mathcal{D})$ (in the case of ensembles, this is equivalent to choosing one among the models)
 2. Propagate forward the learned dynamics according to $x_{t+1} \sim p_{\theta_i}(x_{t+1} | x_t, u_t)$, for all t
 3. Compute (predicted) rewards $\sum_t r(x_t, u_t)$
 4. Repeat steps 1-3 and compute the average reward

$$J(u_1, \dots, u_T) = \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^H r(x_{t,i}, u_t), \text{ where } x_{t+1,i} \sim p_{\theta_i}(x_{t+1,i} | x_{t,i}, u_t)$$



Recap: Model-based RL



Case study: PETS

- Probabilistic Ensembles with Trajectory Sampling
- Key idea:
 - **Model:** Use ensemble of NNs to approximate posterior over model

Deep Reinforcement Learning in a Handful of Trials using Probabilistic Dynamics Models

Kurtland Chua

Roberto Calandra

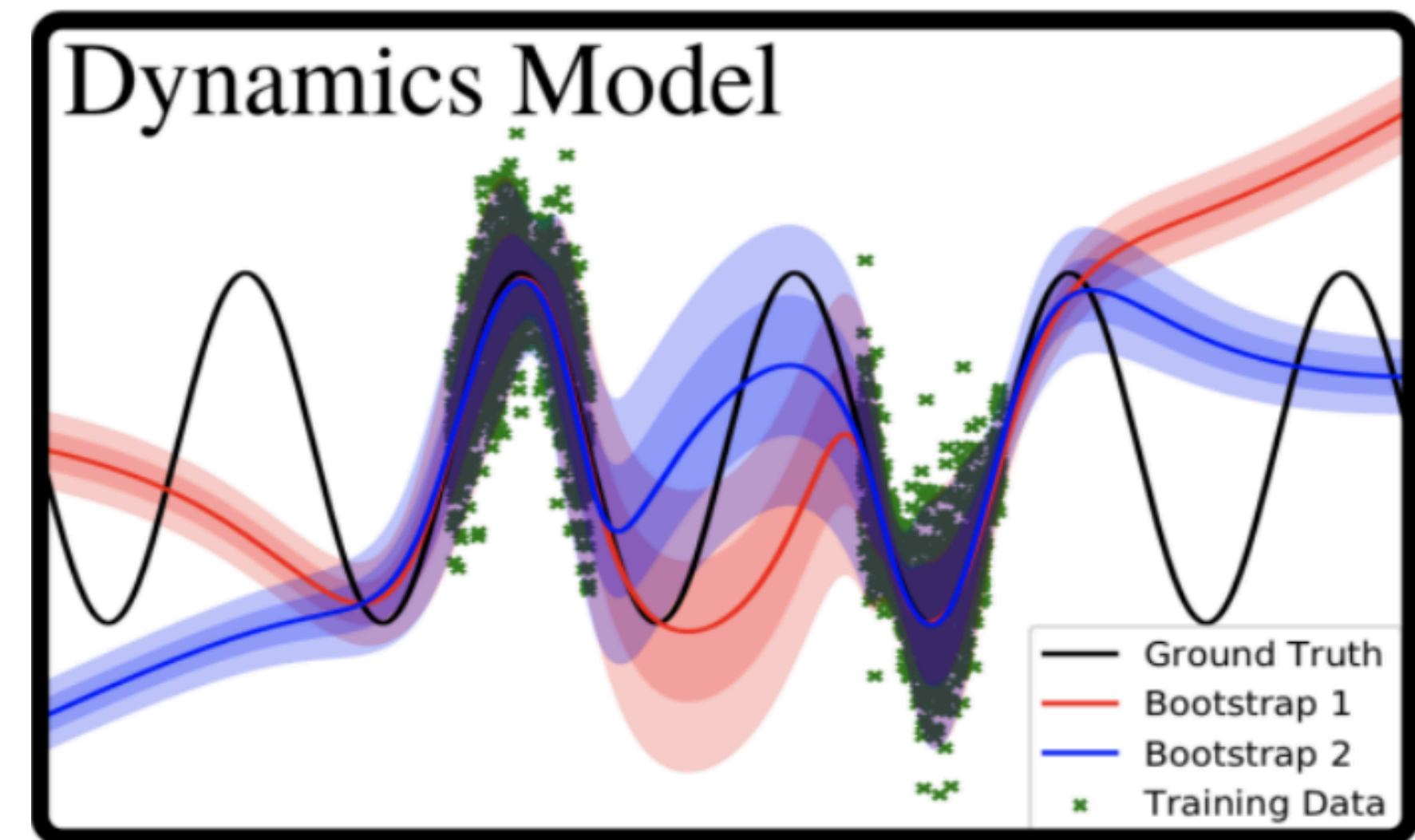
Rowan McAllister

Sergey Levine

Berkeley Artificial Intelligence Research

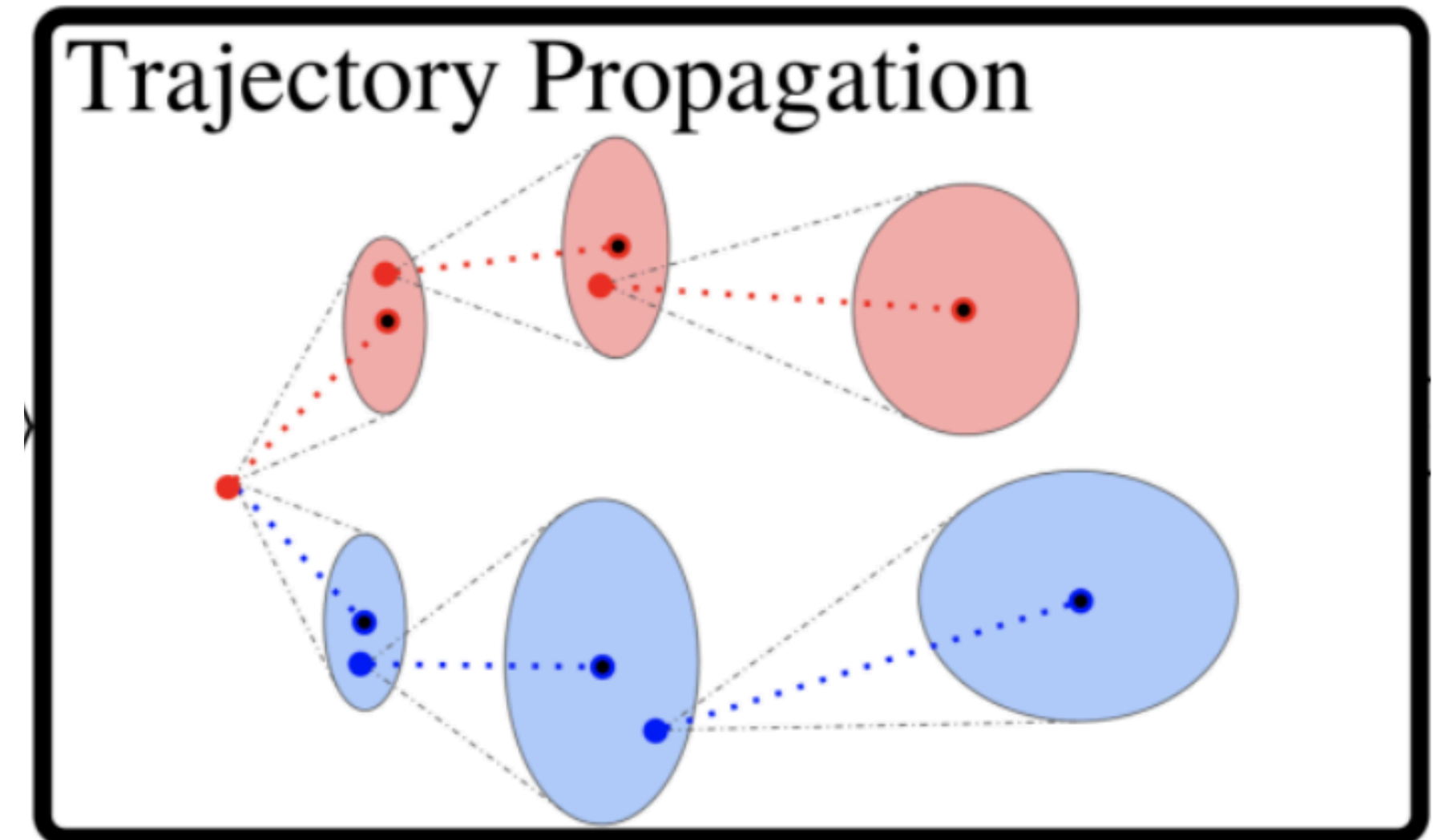
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Case study: PETS

- Probabilistic Ensembles with Trajectory Sampling
- Key idea:
 - **Model:** Use ensemble of NNs to approximate posterior over model
 - **Propagation:** sample different models and use them to generate predictions of different “futures”



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Case study: PETS

- Probabilistic Ensembles with Trajectory Sampling
- Key idea:
 - **Model:** Use ensemble of NNs to approximate posterior over model
 - **Propagation:** sample different models and use them to generate predictions of different “futures”
 - **Planning:** apply MPC (compute action sequence via sampling, i.e., cross-entropy method (CEM))

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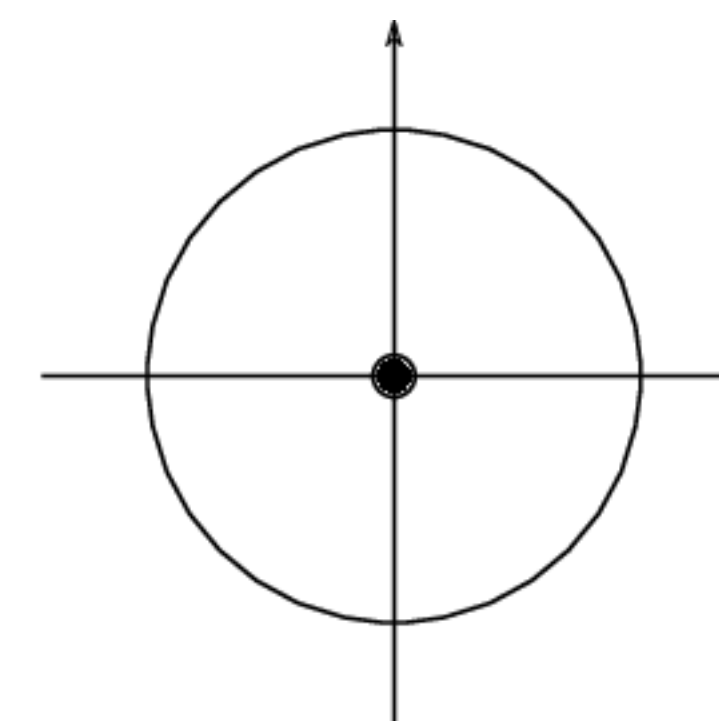
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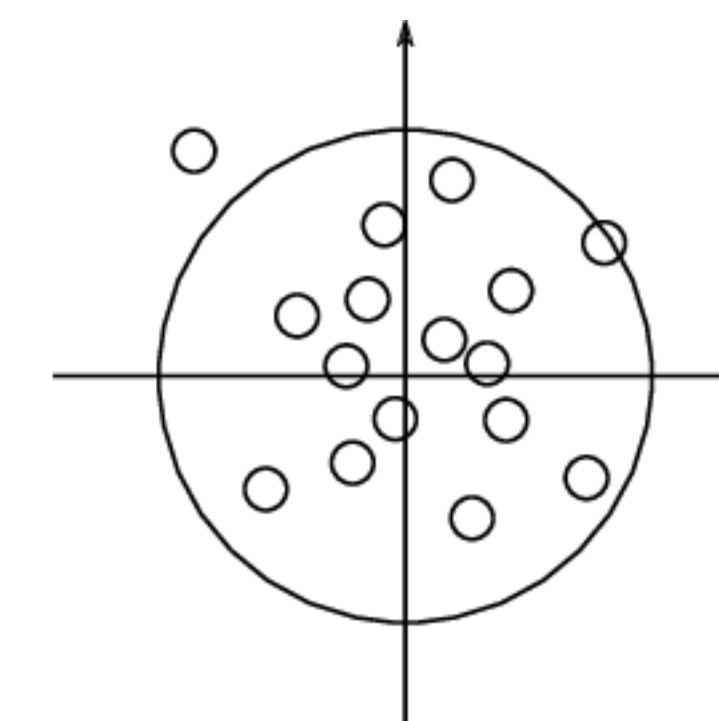
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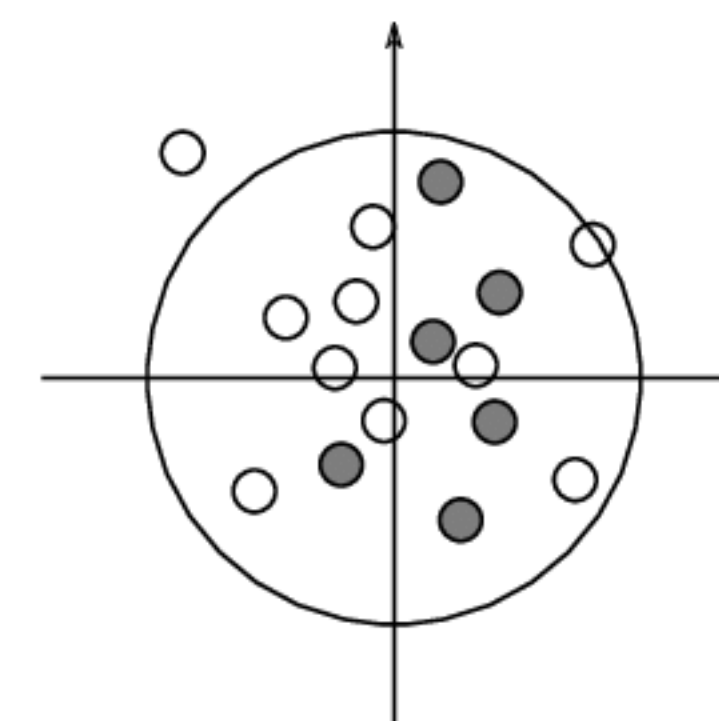
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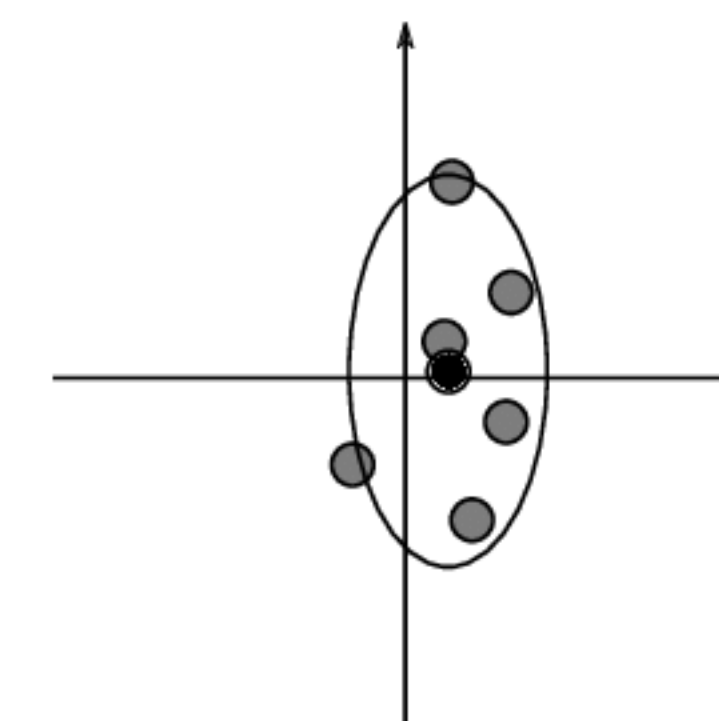
1. Start with the normal distribution $N(\mu, \sigma^2)$



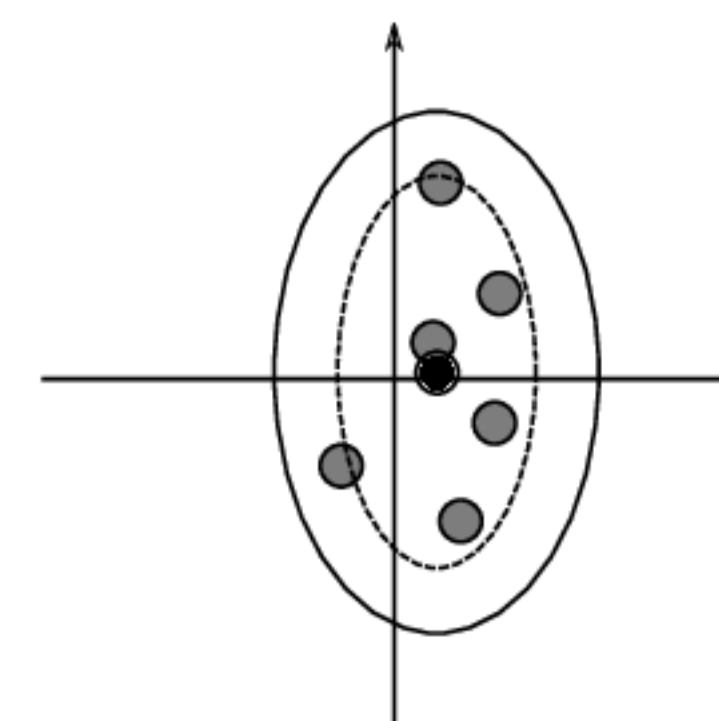
2. Generate N vectors with this distribution



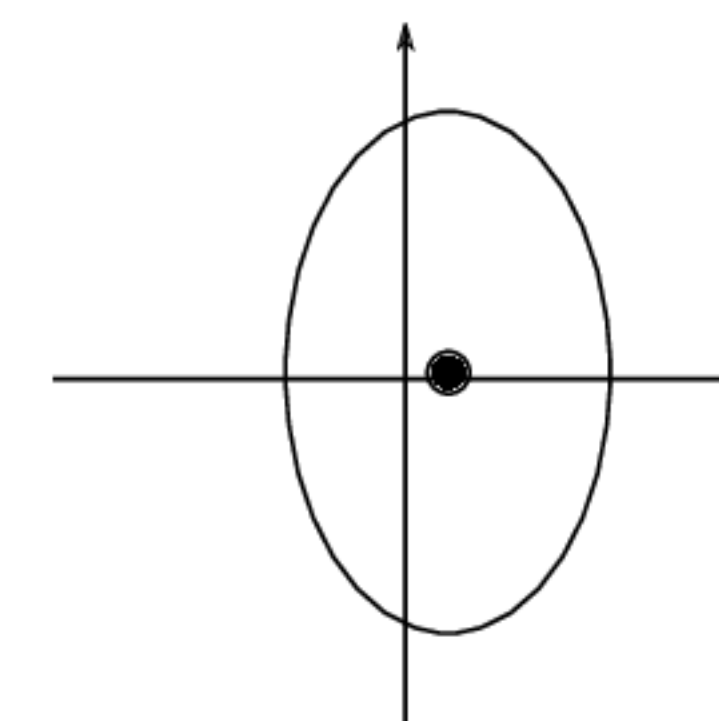
3. Evaluate each vector and select a proportion ρ of the best ones. These vectors are represented in grey



4. Compute the mean and standard deviation of the best vectors

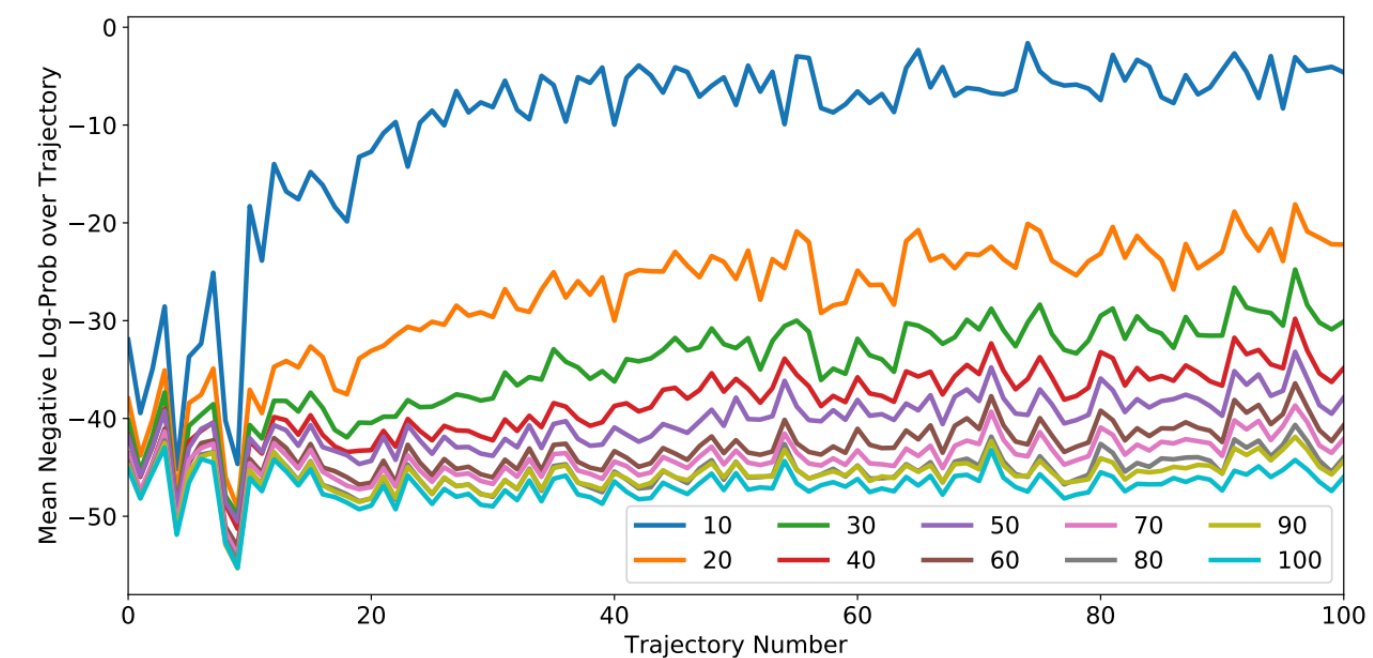
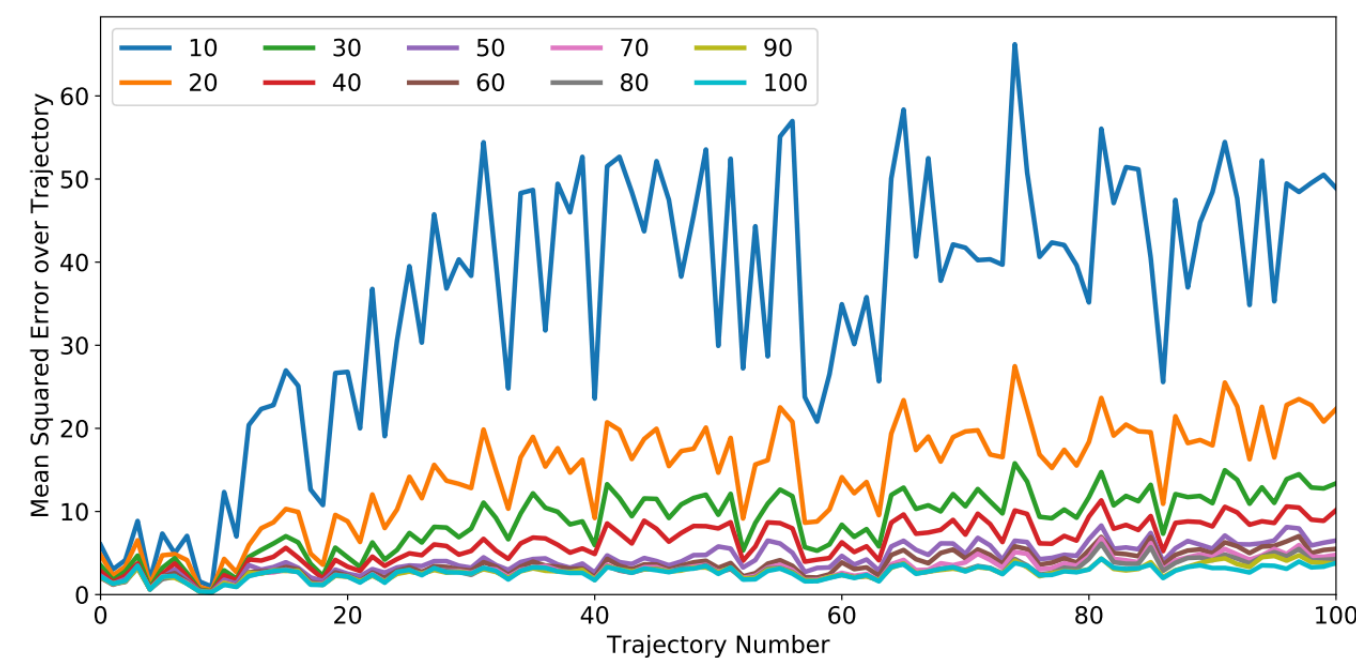
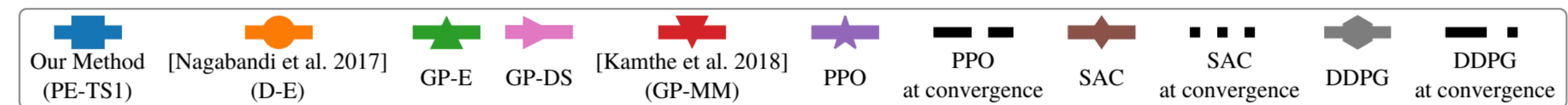
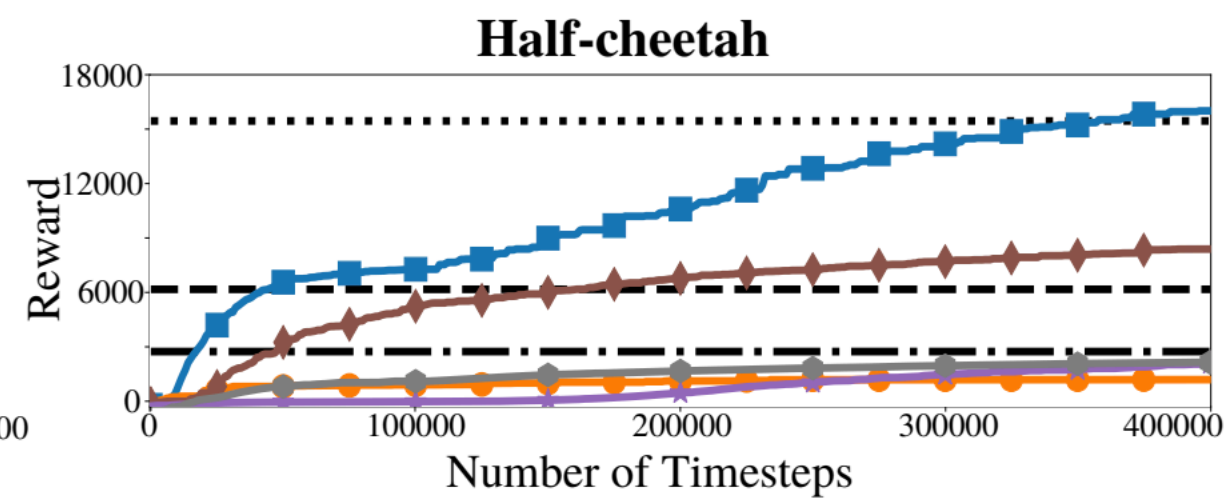
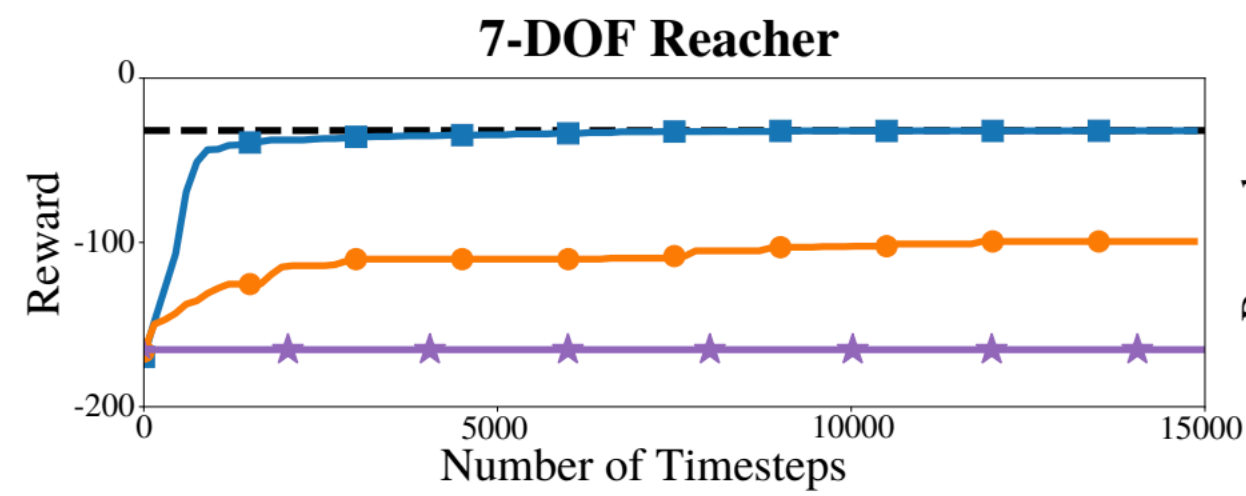
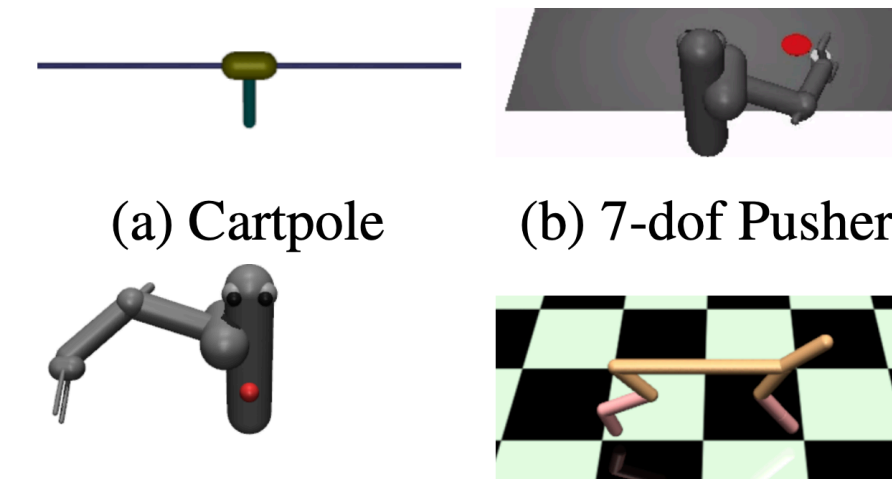
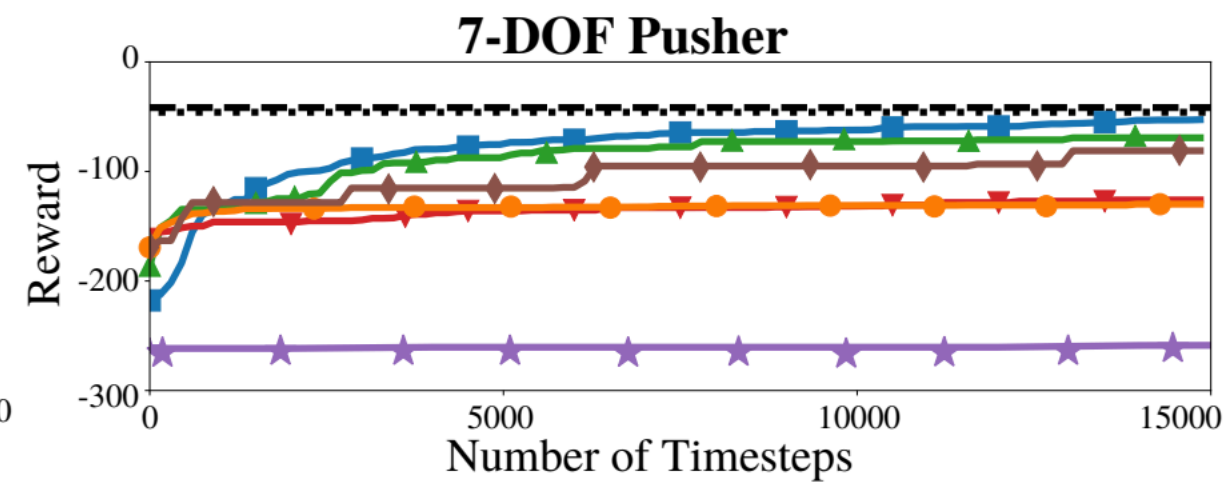
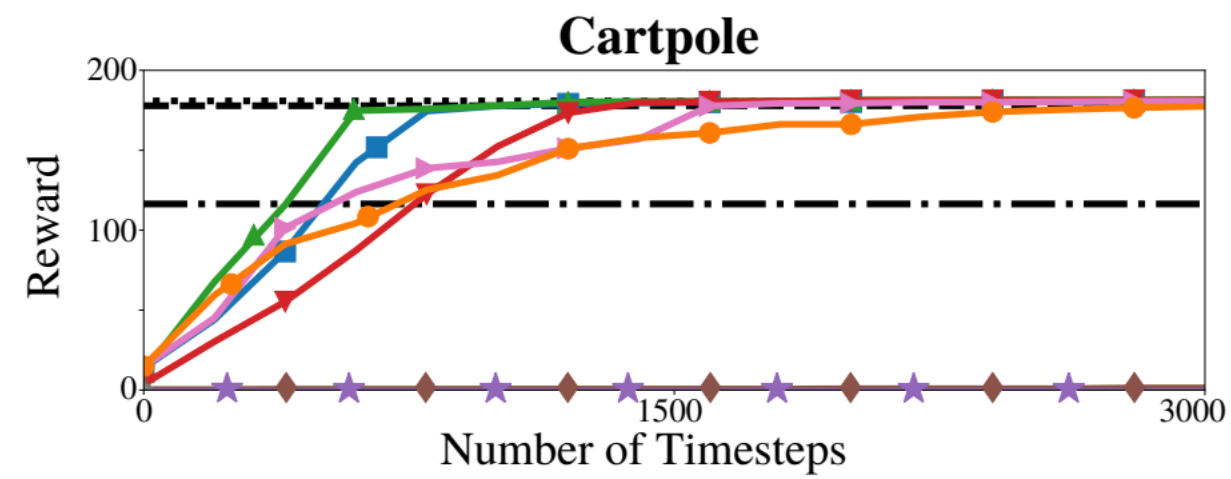


5. Add a noise term to the standard deviation, to avoid premature convergence to a local optimum



6. This mean and standard deviation define the normal distribution of next iteration

Case study: PETS



(a) Mean squared error.

(b) Negative log likelihood.

Why model-based?

- Pros:
 - Sample efficiency
 - Improved multi-task performance
 - Transitions provide strong supervision (opposed to e.g., sparse reward)
- Cons:
 - Optimize the wrong objective
 - Can converge to worse performance if model is wrong
 - Can be difficult to train with high-dimensional states/observations (e.g., images)

Outline

Last week

Approach 1:

“Learn a model and use it to plan”

General idea

- Integrating planning and learning

“Dyna-style” algorithms

- Dyna-Q & Extensions

Remarks

Approach 2:

“Learn a model and improve model-free learning”

A bird's eye view of previous lectures

- Value-based methods: learn **value functions** from experience
- Policy optimization: learn **policies** from experience
- Previous lecture: learn a **model** from experience (and **plan** to construct a policy)
- Today: integrate learning and planning into a single architecture

Note:

- Last week we used the term model to describe a dynamics model, i.e.,

$$x_{t+1} \sim p_{\theta}(x_{t+1} | x_t, u_t)$$

- In general, we can assume the model to represent the *unknowns* in our MDP $\mathcal{M} = (X, U, P, R)$

$$x_{t+1} \sim p_{\theta}(x_{t+1} | x_t, u_t)$$

$$R_t = r_{\theta}(x_t, u_t)$$

Examples of models:

- Table look-up
- Linear
- GP
- Neural network, ...

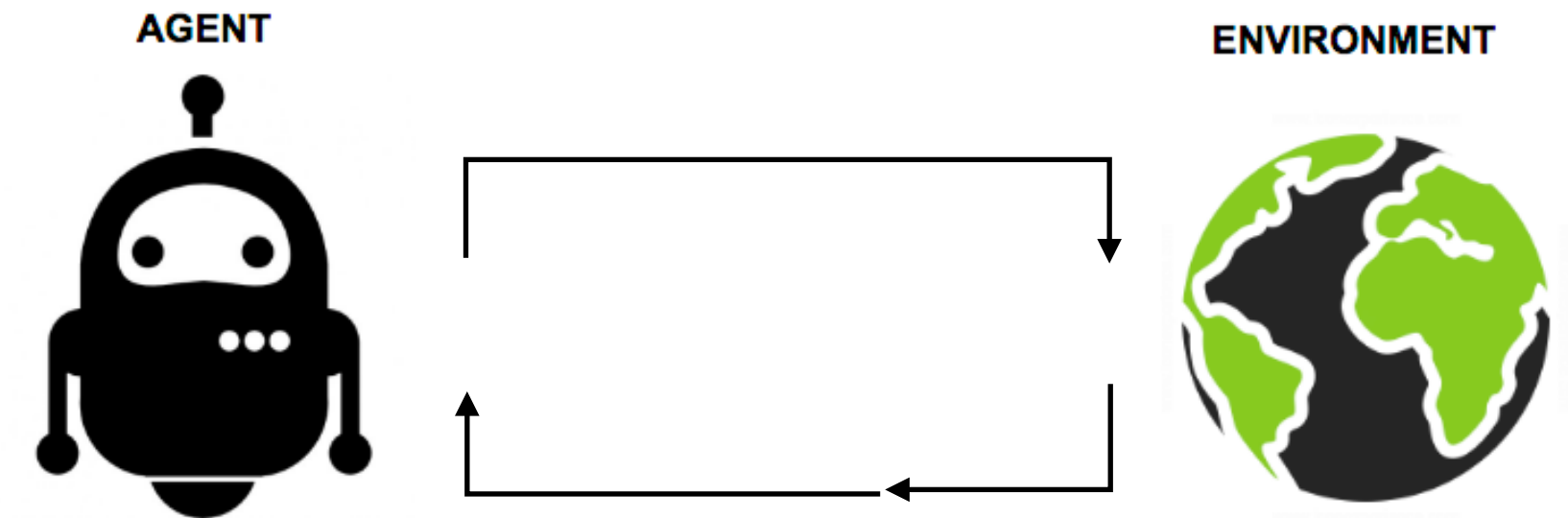
Different sources of experience

- Having a model enables us to consider two sources of experience

Real experience: sampled from the environment (true MDP)

- Interacting with the environment provides us with samples from the true MDP $\mathcal{M} = (X, U, P, R)$

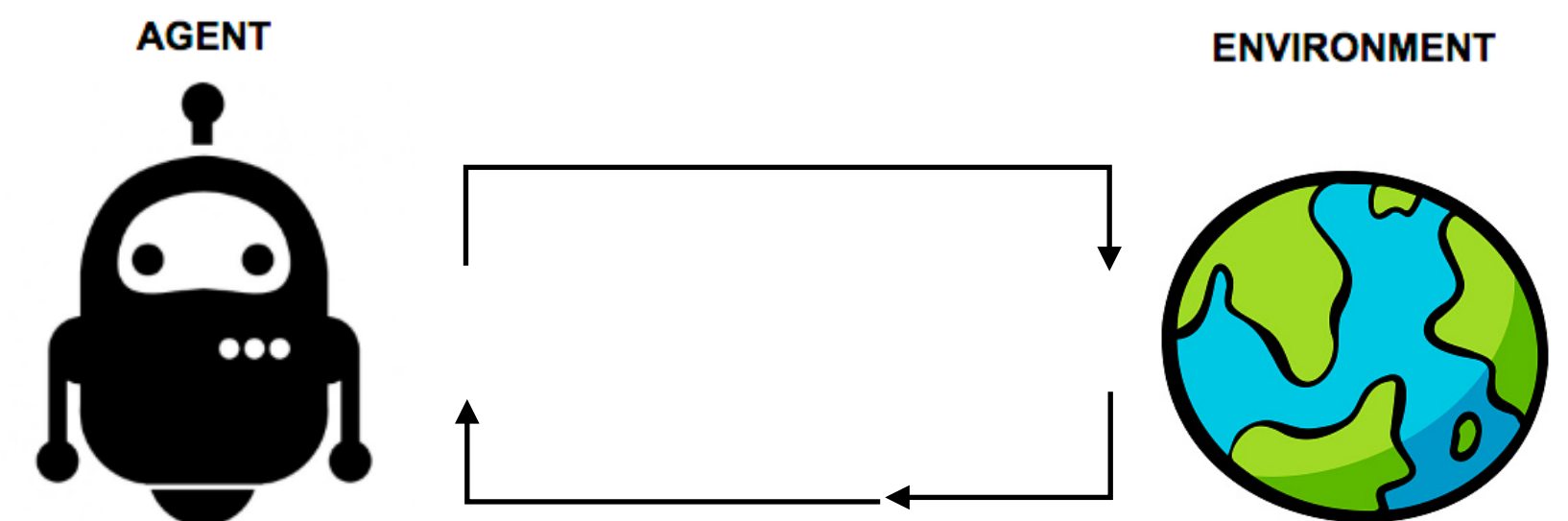
$$x_{t+1} \sim P(x_{t+1} | x_t, u_t)$$
$$R_t = R(x_t, u_t)$$



Simulated experience: sampled from the model (approximate MDP)

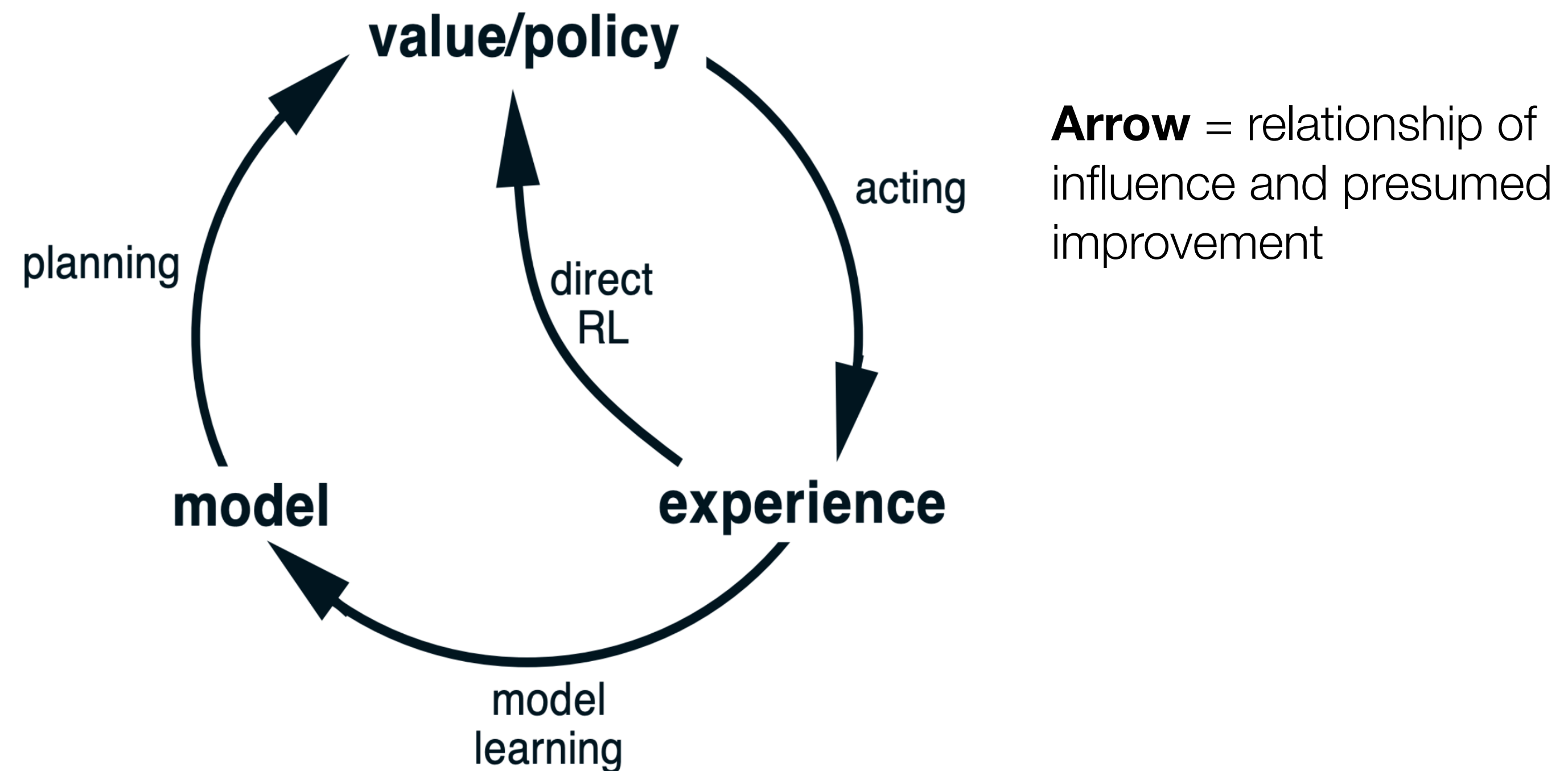
- Simulating transitions through the model provides us with samples from an approximation of the MDP

$$x_{t+1} \sim p_{\theta}(x_{t+1} | x_t, u_t)$$
$$R_t = r_{\theta}(x_t, u_t)$$



A general recipe for model-based acceleration

1. Interact with the environment to generate a dataset of transitions $\mathcal{D} = \{(x_t, u_t, r_t, x_{t+1})\}$
2. Fit dynamics/reward model to \mathcal{D}
3. Generate simulated experience under your model and use model-free algorithms



Example: MBRL via policy gradient

- In PO, we defined the policy gradient via (variations of) the following equation:

$$\nabla_{\theta} J(\theta) \approx \frac{1}{N} \sum_{i=1}^N \sum_{t=1}^T \nabla_{\theta} \log \pi_{\theta}(u_{i,t} | x_{i,t}) Q_{\phi}(x_{i,t}, u_{i,t})$$

where used real experience (in the form of trajectories of interactions with the environment) to practically approximate the expectations

- We could consider the following scheme:

1. Run base policy $\pi_0(u_t | x_t)$ in the environment (e.g., random policy, exploration policy) and collect dataset of transitions

$$\mathcal{D} = \{(x_t, u_t, x_{t+1})_i\}$$

2. Fit dynamics model to data to minimize error (or equivalently, maximize (log) likelihood)

$$\theta^* = \arg \min_{\theta} \sum_i \left\| f_{\theta}(x_t, u_t) - x_{t+1} \right\|^2$$

3. Use the learned model to generate simulated trajectories $\{\tau_i\}$ through policy π_{θ}
4. Use $\{\tau_i\}$ to improve π_{θ} via policy gradient

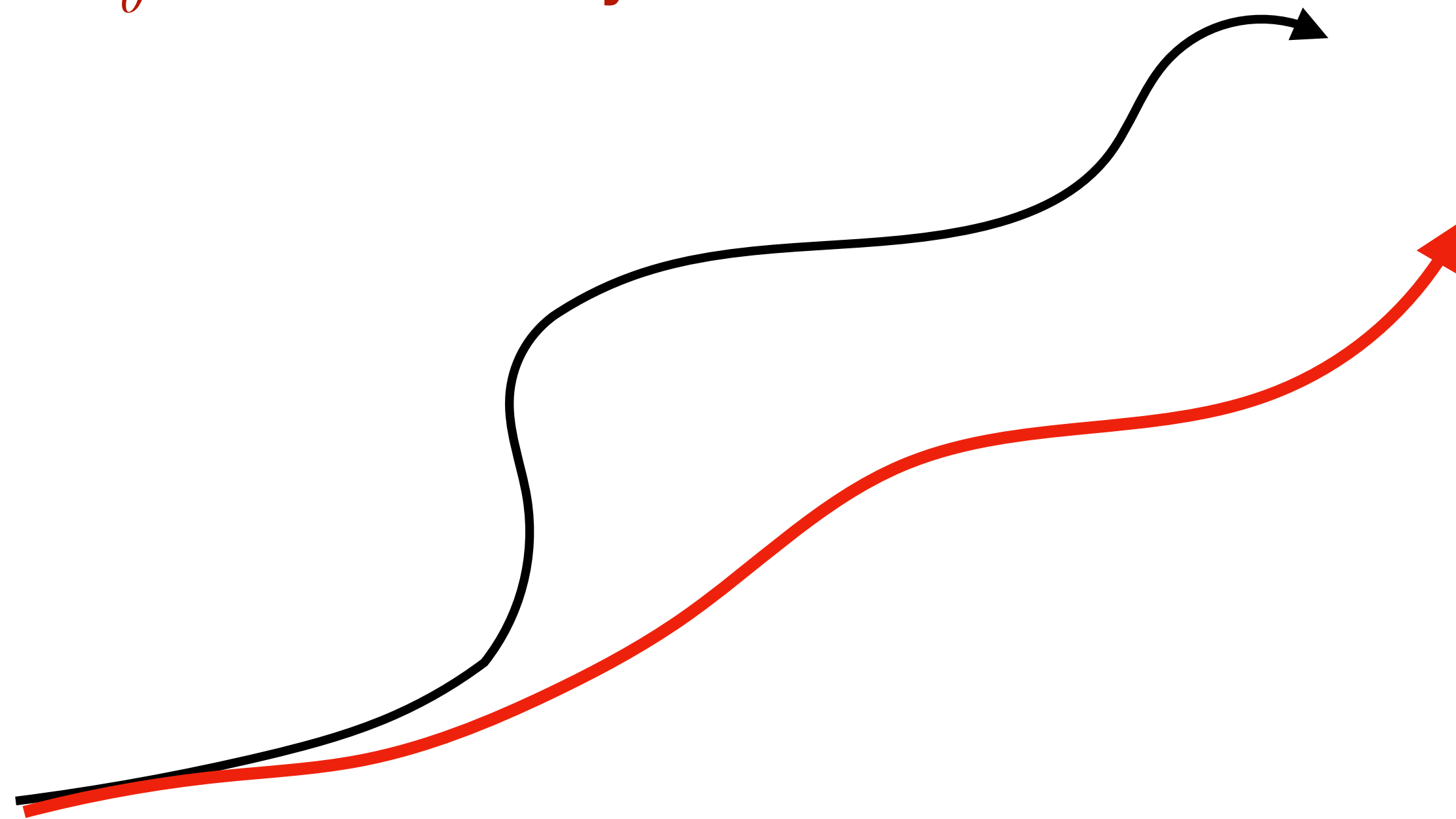
Question

What is a potential problem with this?

Issue with (long) model-based rollouts

Run π_θ with real dynamics

Run π_θ with estimated dynamics



- We want to avoid long model-based rollouts, as these will necessarily incur in accumulating error
- At the same time, short rollouts do not guarantee exploration of “later timesteps”

Dyna-Q

Initialize $Q(x, u)$ and $Model(x, u)$, $\forall x \in X, \forall u \in U$,

Repeat (for each episode):

(a) $x_t \leftarrow$ current (non-terminal) state

(b) Choose u_t from x_t using policy derived from Q (e.g., ϵ -greedy)

(c) Take action u_t , observe r_t, x_{t+1}

(d) $Q(x_t, u_t) \leftarrow Q(x_t, u_t) + \alpha \left(r_t + \gamma \max_{u'_{t+1}} Q(x_{t+1}, u'_{t+1}) - Q(x_t, u_t) \right)$

(e) $Model(x_t, u_t) \leftarrow x_{t+1}, r_t$

(f) Repeat n times:

$x_t \leftarrow$ sample random state previously observed

$u_t \leftarrow$ sample random action previously taken in x_t

$x_{t+1}, r_t \leftarrow Model(x_t, u_t)$; predict through model

$Q(x_t, u_t) \leftarrow Q(x_t, u_t) + \alpha \left(r_t + \gamma \max_{u'_{t+1}} Q(x_{t+1}, u'_{t+1}) - Q(x_t, u_t) \right)$

until x_t is terminal

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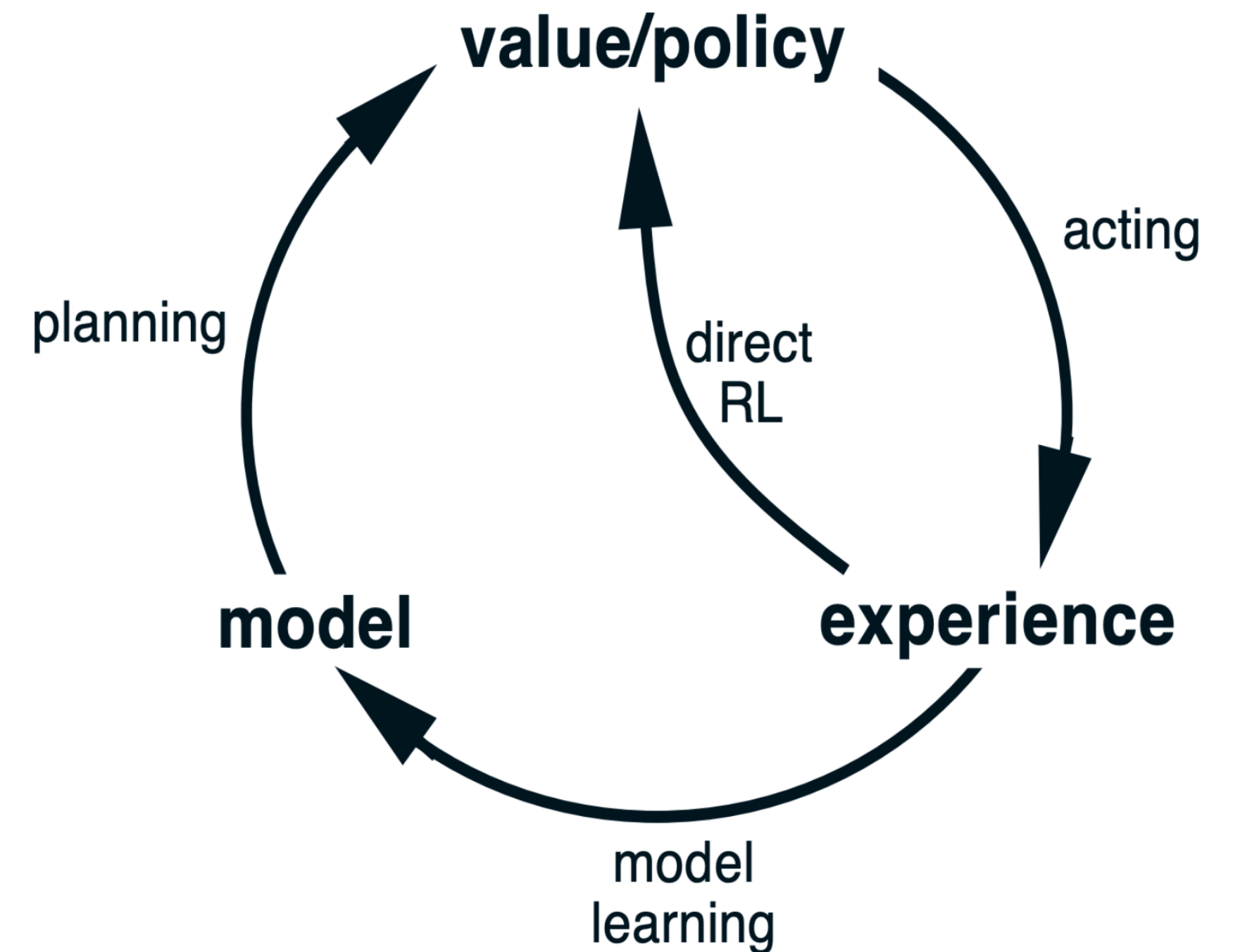
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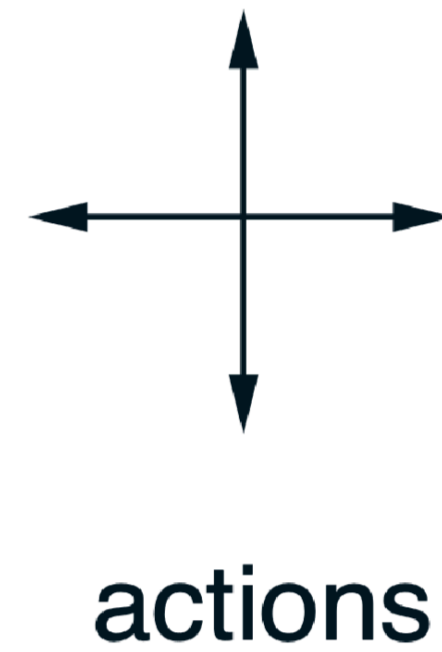
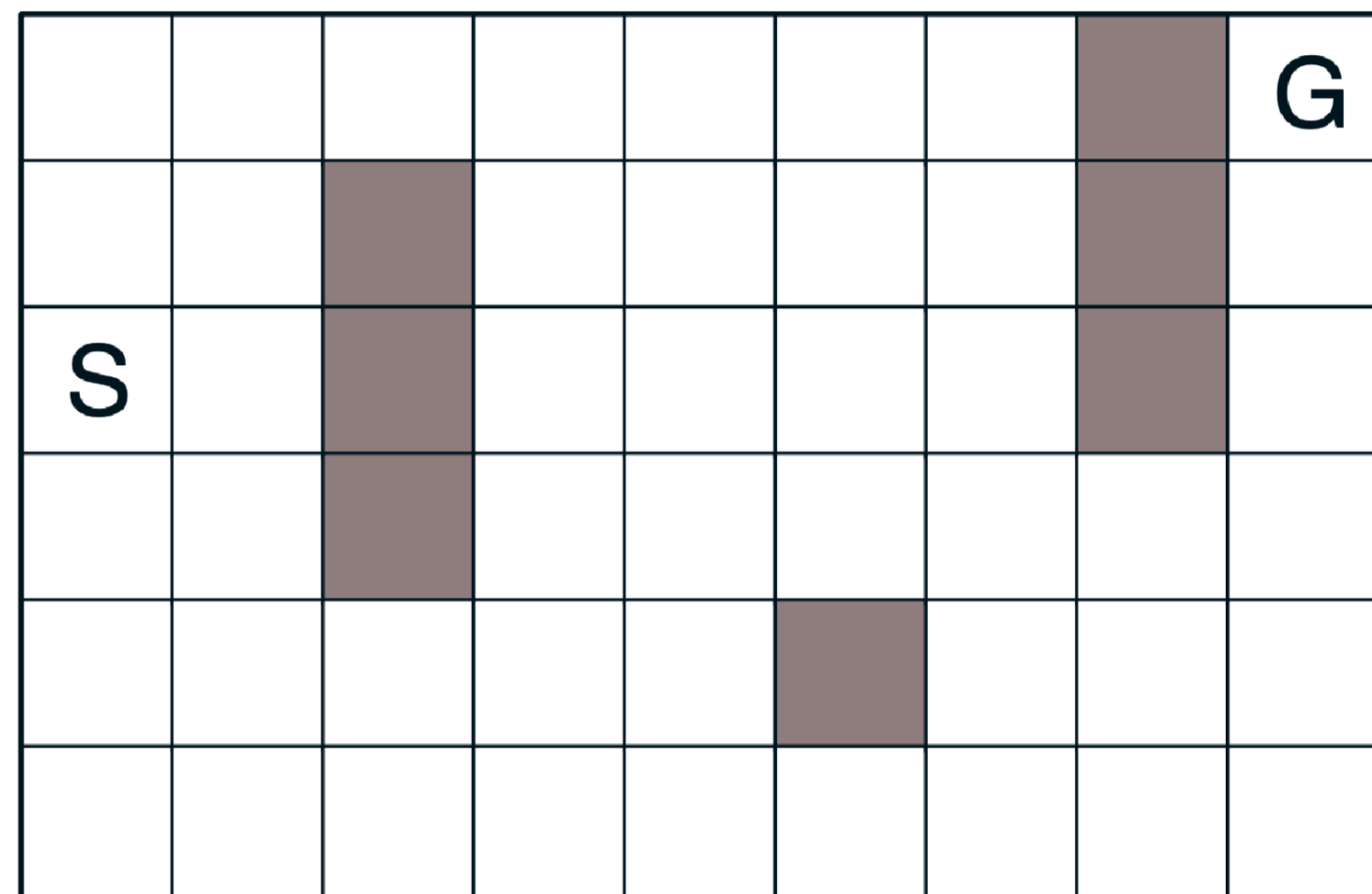
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until x_t is terminal



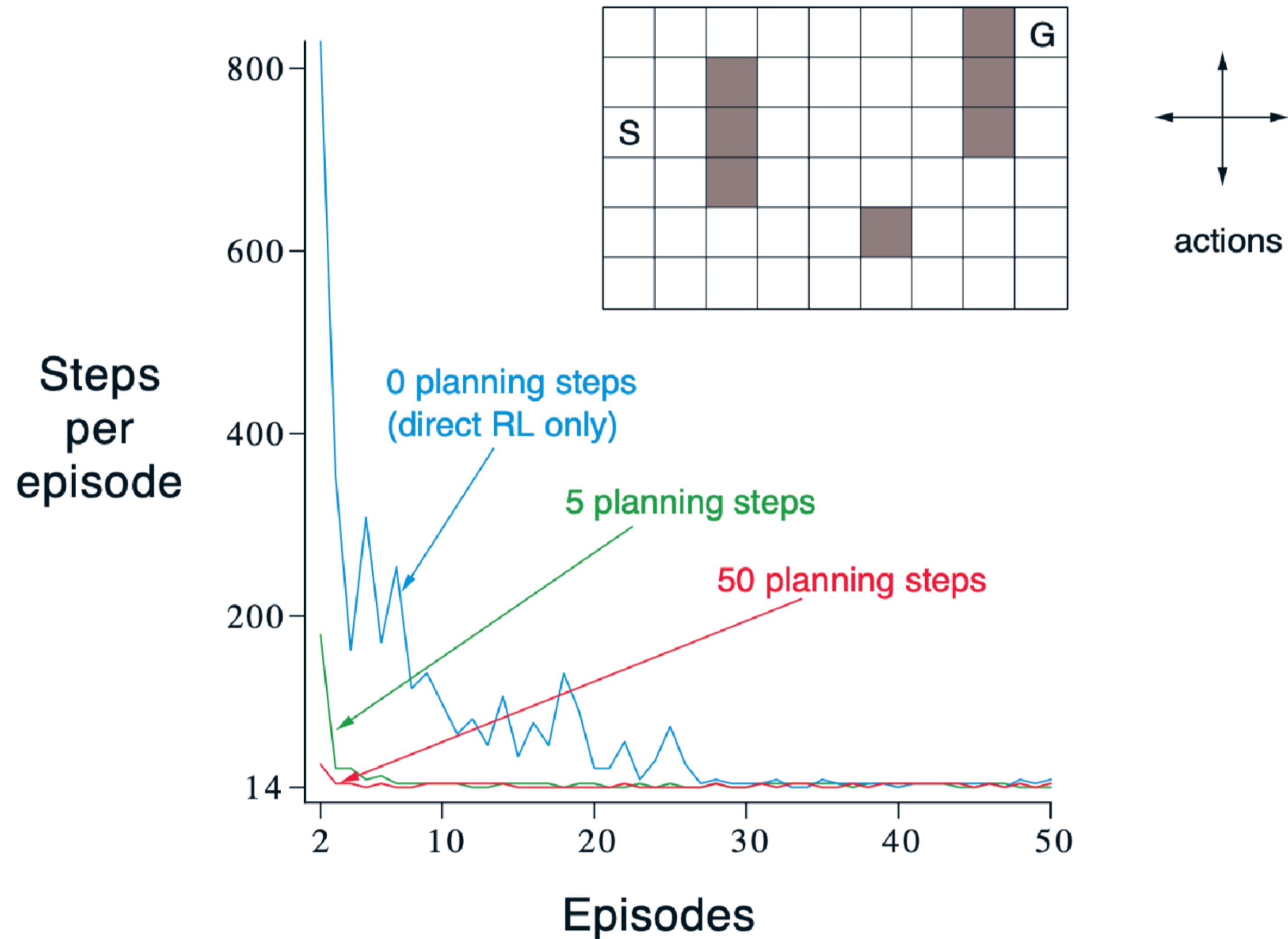
- **Model** = Tabular model
- **Direct RL** = Q-learning
- **Planning** = 1-step

Example: Dyna-maze



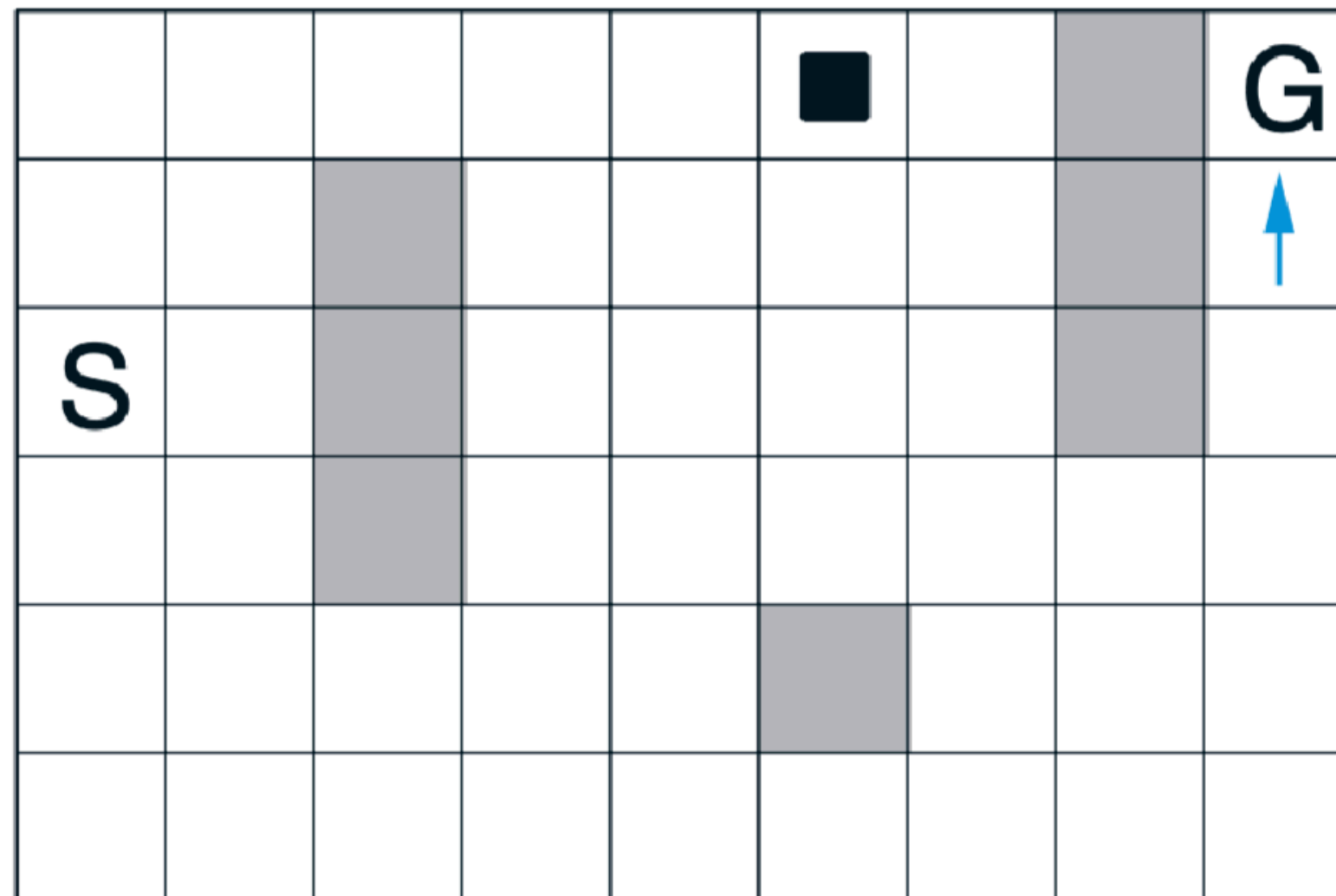
- 47 states, 4 actions
- Deterministic dynamics
- Reward = 0 everywhere, except +1 on G
- $\gamma = 0.95$
- Zero-initialized Q, $\alpha = 0.1, \epsilon = 0.1$

Example: Dyna-maze

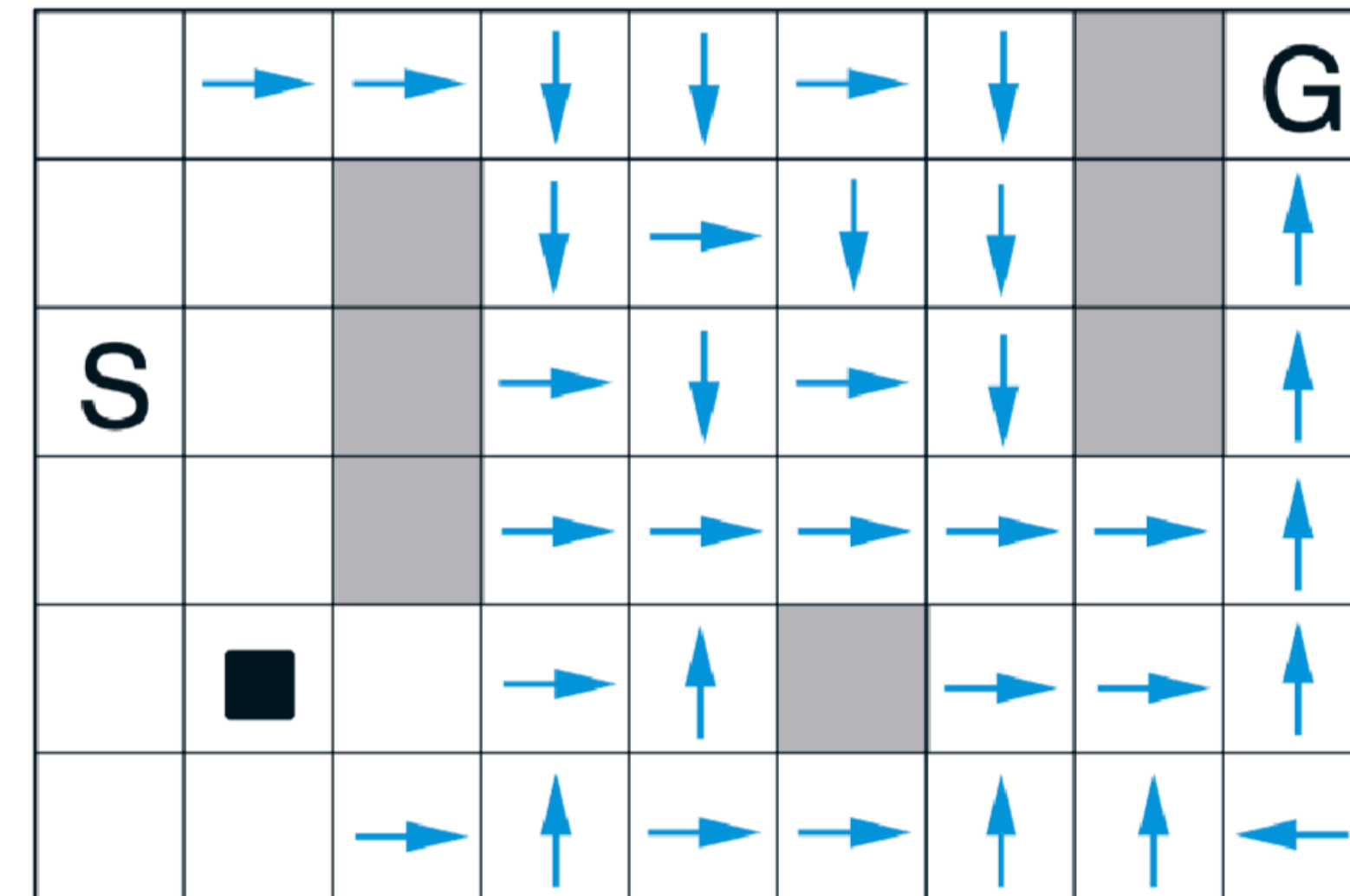


Example: Dyna-maze

WITHOUT PLANNING ($n=0$)



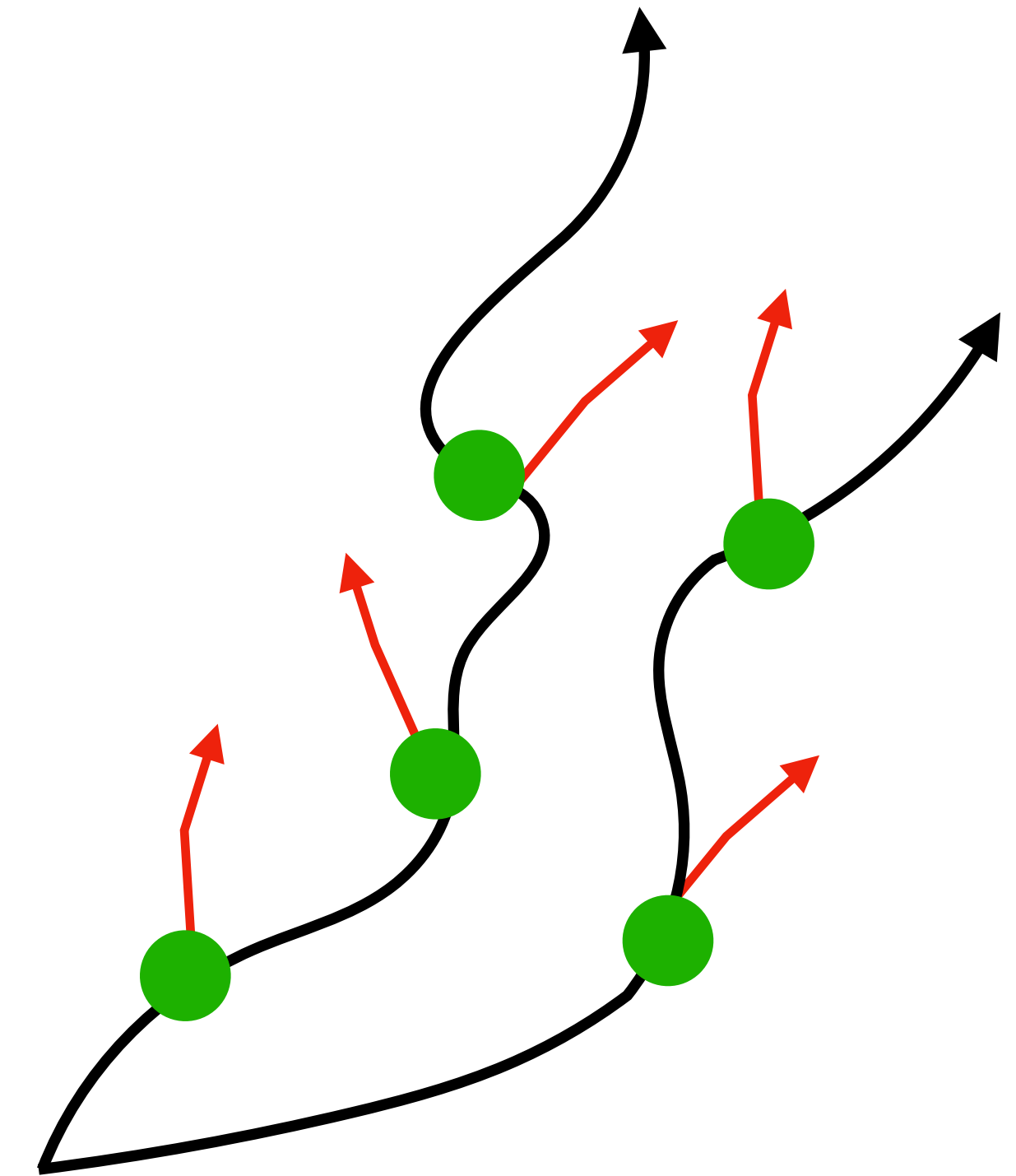
WITH PLANNING ($n=50$)



- The plot compares the policies found by Dyna-Q with and without planning, half-way through the second episode
- Without planning ($n = 0$), each episode adds only one additional step to the policy, and so only one step (the last) has been learned so far
- With planning, again only one step is learned during the first episode, but during the second episode planning allows to develop an extensive policy that will reach almost back to the start state

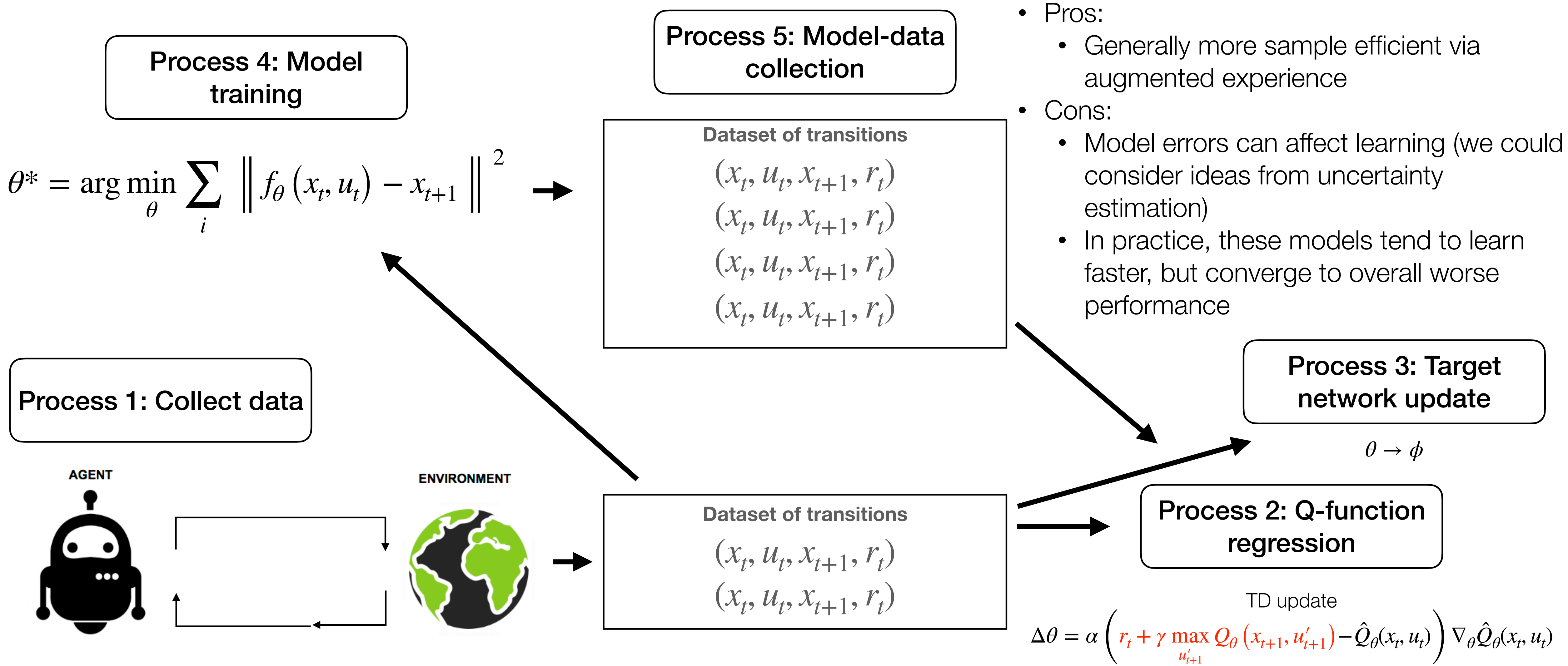
“Dyna-style” algorithms

- Dyna-Q represents a specific choice of model, planning, direct RL algorithm, etc.
- More generally, we can define the following recipe for “dyna-style” algorithms
 1. Collect data $\{(x_t, u_t, r_t, x_{t+1})\}$
 2. Learn dynamics / reward model, i.e., $p_\theta(x_{t+1} | x_t, u_t), r_\theta(x_t, u_t)$
 3. Repeat n times
 1. Sample x_t from buffer
 2. Choose action u_t (from dataset, π , random, exploration, etc.)
 3. Simulate dynamics / reward $\hat{x}_{t+1} \sim p_\theta(x_{t+1} | x_t, u_t), \hat{r}_t = r_\theta(x_t, u_t)$
 4. Train on $\{(x_t, u_t, \hat{r}_t, \hat{x}_{t+1})\}$ via model-free RL
 5. Optionally, take k more model-based steps



- Only uses short roll-outs
- While observing diverse states

Example: model-based acceleration of DQN



- Pros:
 - Generally more sample efficient via augmented experience
- Cons:
 - Model errors can affect learning (we could consider ideas from uncertainty estimation)
 - In practice, these models tend to learn faster, but converge to overall worse performance

Case study: PILCO

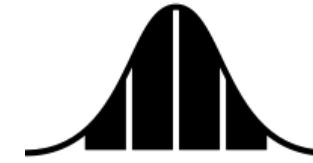
- Deisenroth and Rasmussen, *Probabilistic inference for learning control*, ICML 2011
- Approach: use Gaussian process for dynamics model
 - Gives measure of *epistemic* uncertainty
 - Extremely sample efficient
- Pair with arbitrary (possibly nonlinear) policy
- By propagating the uncertainty in the transitions, capture the effect of small amount of data



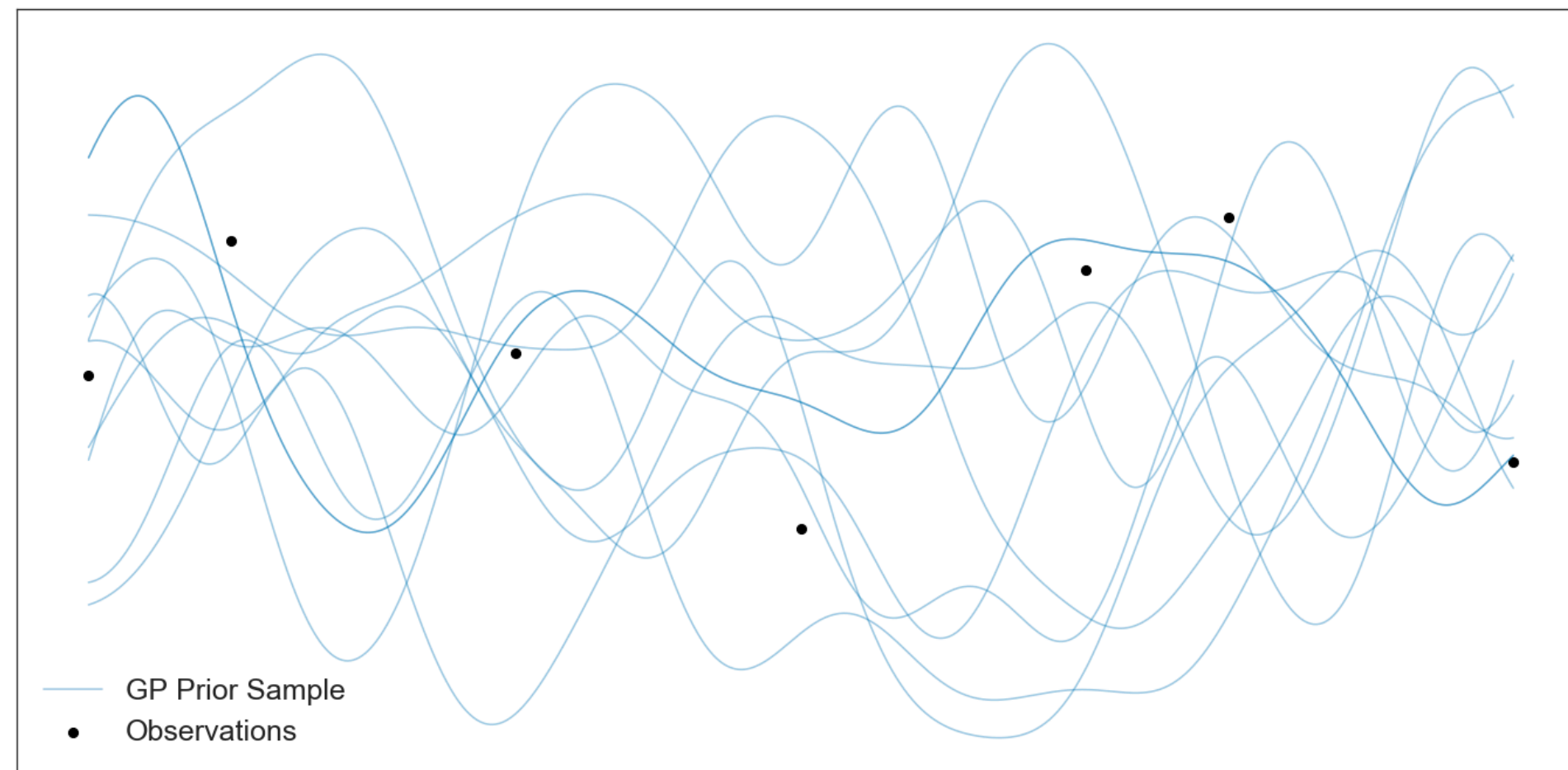
Gaussian process reminder

- Represent “distribution over functions”

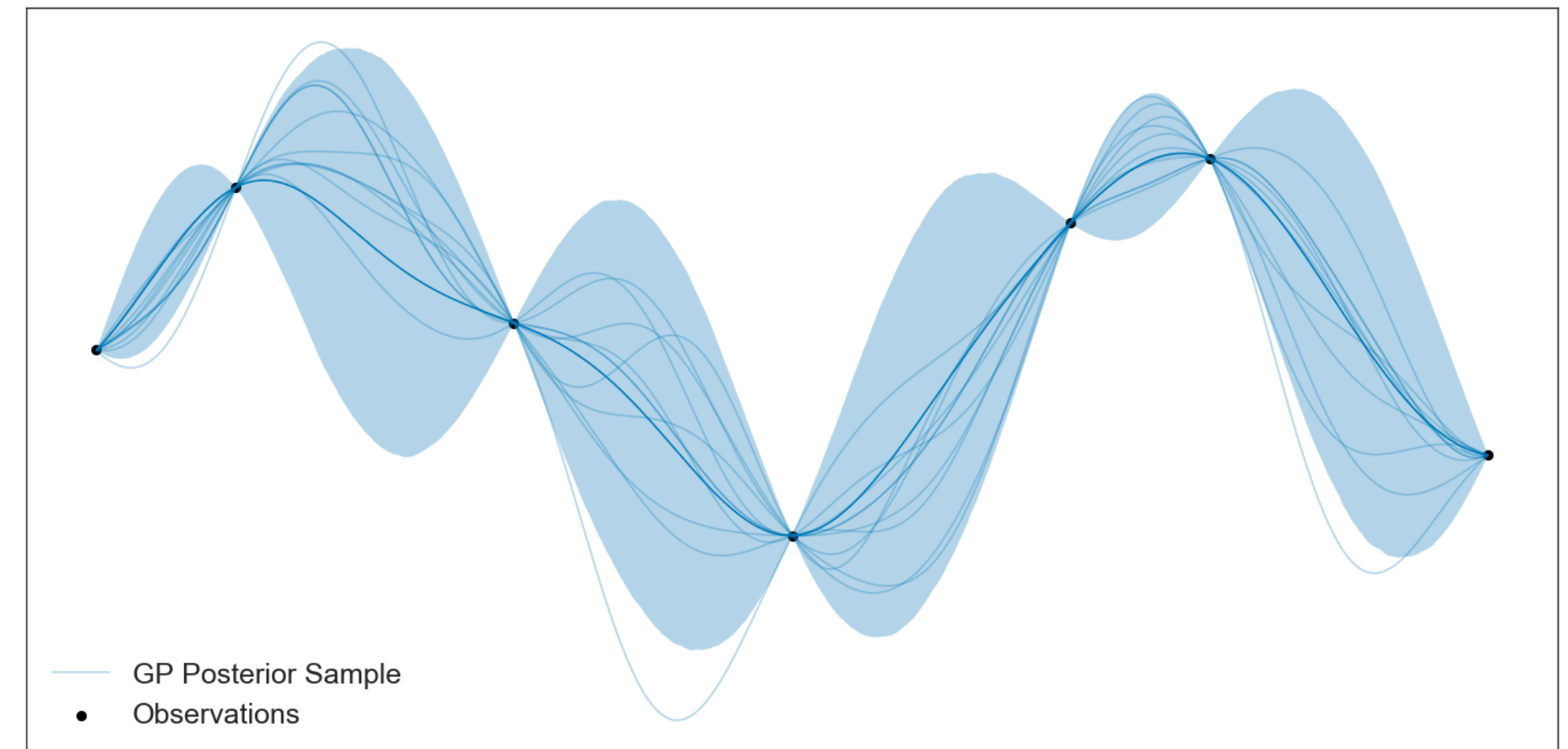
Bayesian inference



Samples from **prior** distribution

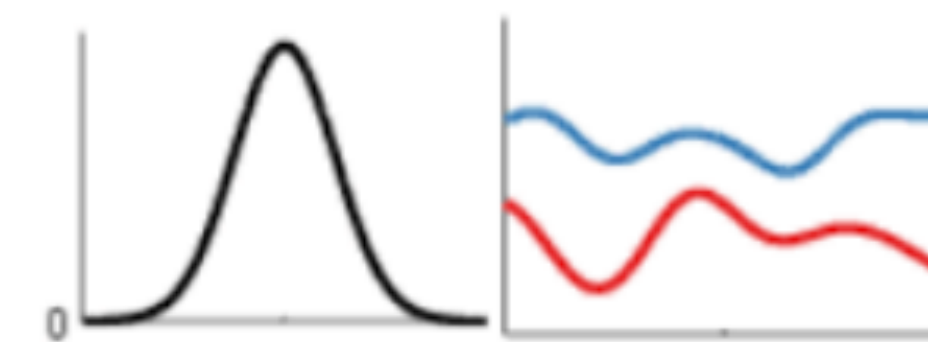


Samples from **posterior** distribution



- Typically, initialize with zero mean; behavior determined entirely by kernel
- Standard kernel choice: squared exponential, used in PILCO
 - Has smooth interpolating behavior

Squared Exponential Kernel

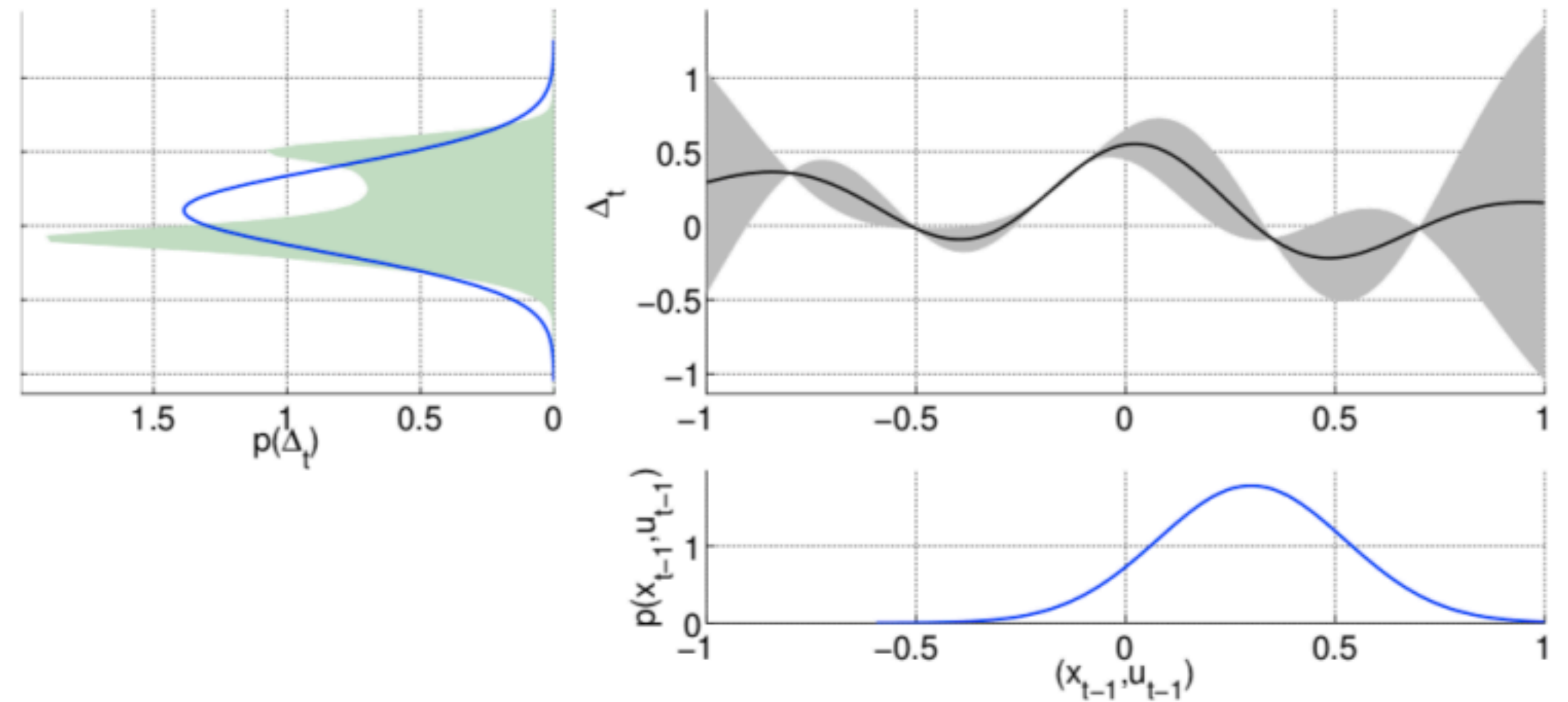


A.K.A. the Radial Basis Function kernel,

$$k_{SE}(x, x') = \sigma^2 \exp\left(-\frac{(x-x')^2}{2\ell^2}\right)$$

Case study: PILCO

- For GP conditioned on data, one step prediction is Gaussian
- But, need to make multistep predictions: so, need to derive multi-step predictive distribution
- Turn to approximating distribution at each time with a Gaussian via moment matching



Case study: PILCO

- All algorithm design choices made to ensure analytical tractability:
- Because of the squared exponential kernel, mean and variance can be computed in closed form
- Choose cost:

$$c(\mathbf{x}) = 1 - \exp\left(-\|\mathbf{x} - \mathbf{x}_{\text{target}}\|^2 / \sigma_c^2\right)$$

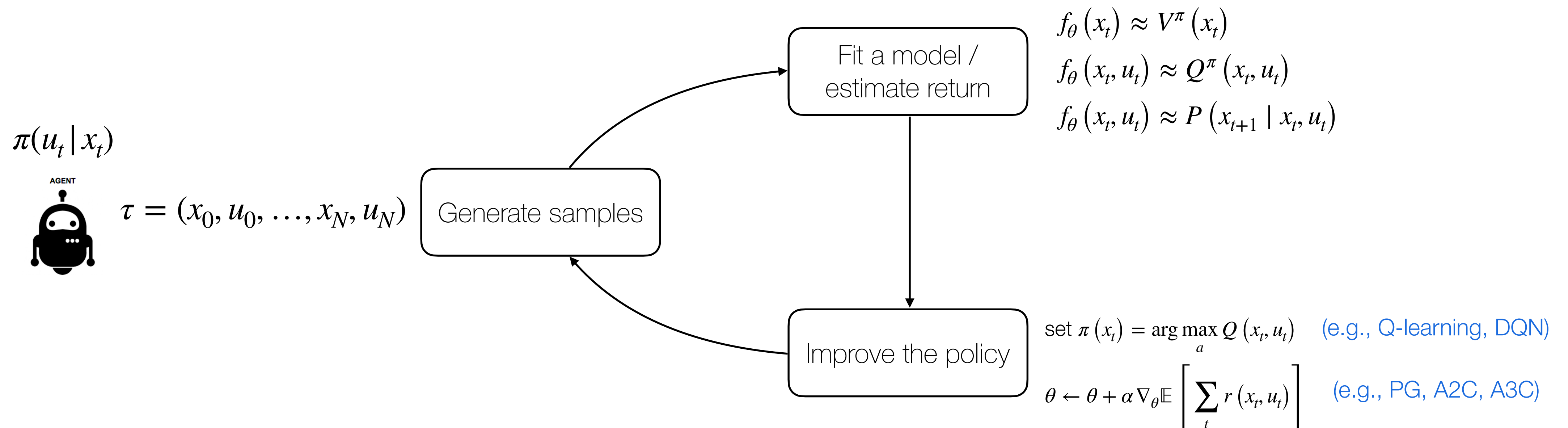
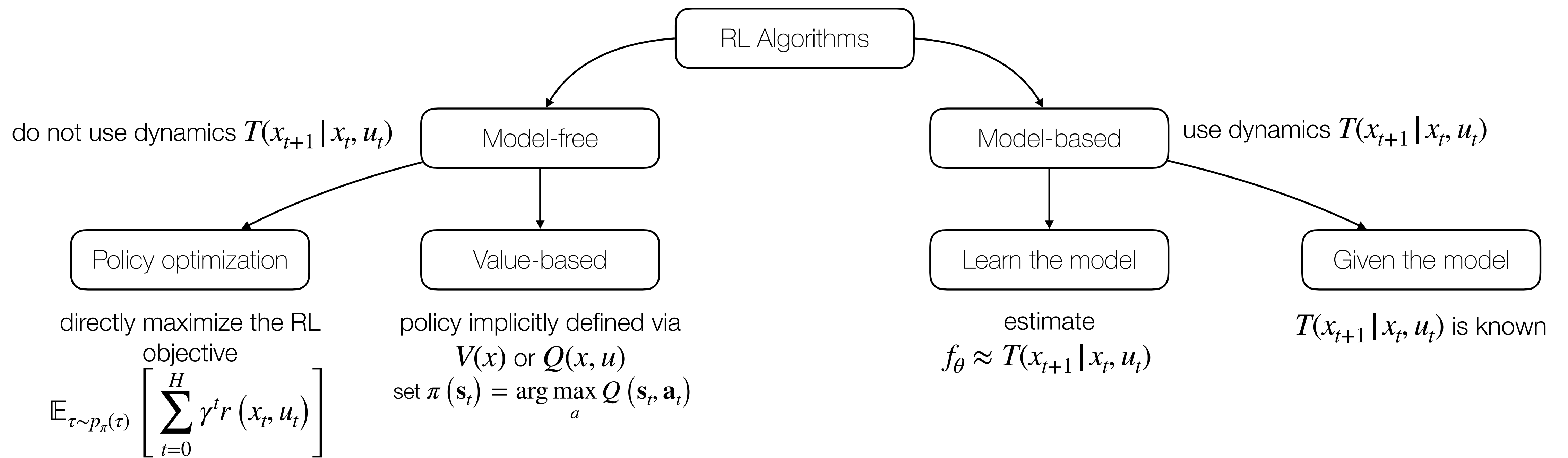
- which is similarly squared exponential; thus expected cost can be computed exactly, factoring in uncertainty
- Choose also radial basis function or linear policy, to enable analytical uncertainty propagation

PILCO (at a high level)

- Uncertainty prop: leverage specific functional forms to derive analytical expressions for mean and variance of trajectory under policy.
- Can use chain rule (aka backprop through time) to compute the gradient of expected total cost w.r.t. policy parameters
- Algorithm:
 - Roll out policy to get new measurements; update model
 - Compute (locally) optimal policy via gradient descent
 - This policy is “local” in the sense of the data we’ve given it, i.e., it’s tailored to the regions of state space it’s seen before; this is more general than “local” in the sense of linearization
 - Repeat

Combining model and policy learning

- We discussed two possible solutions, but there are infinitely many more!
- Very busy research direction! Many topics not covered here
 - Many possible combinations of planning/control, policies, values, and models
- Quite practical: model learning is data efficient and parameterized policy is cheap to evaluate at run time



Next time

- Course recap
- Research presentations