

# Faster RL by Freezing Slow States

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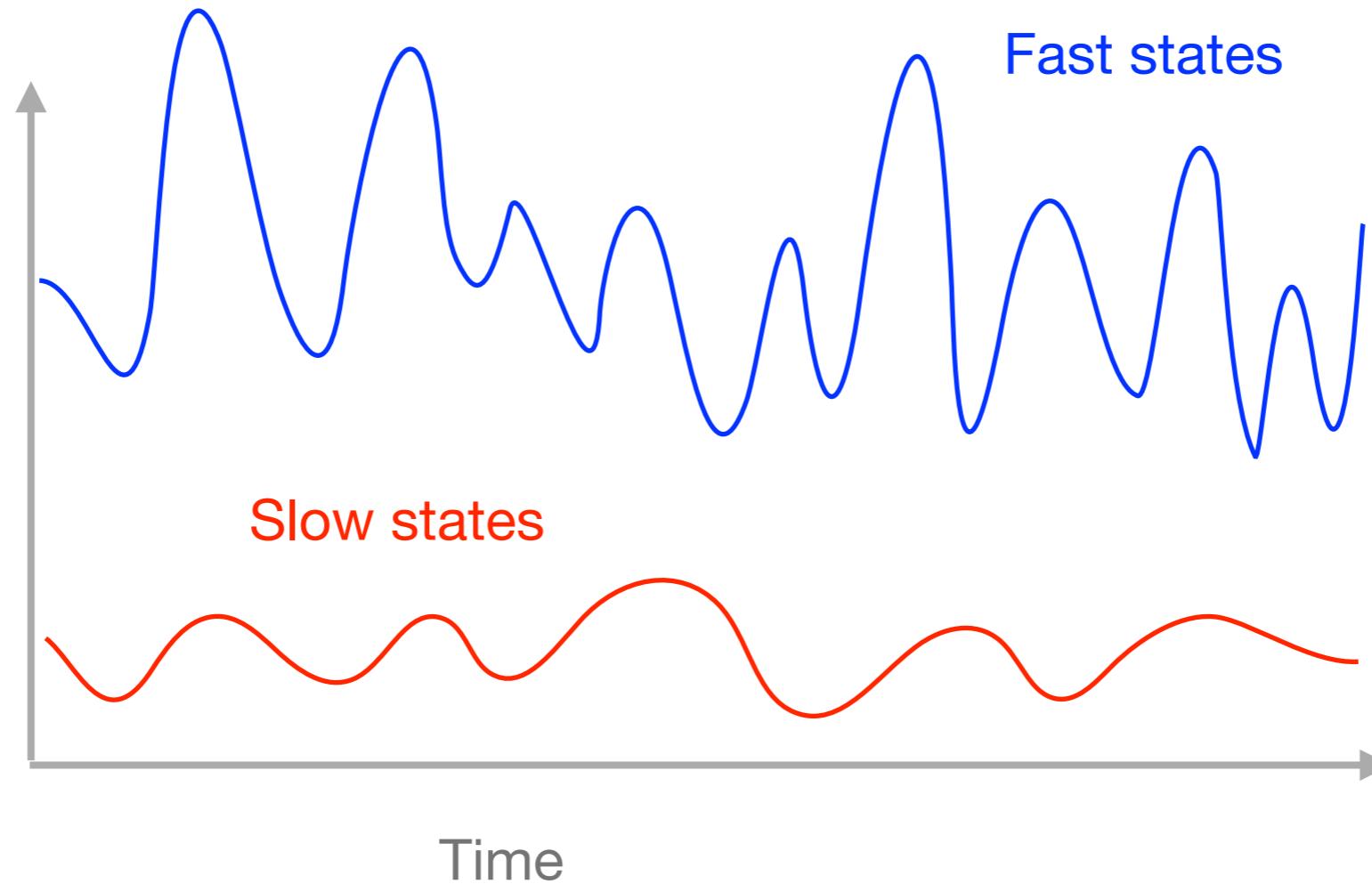
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*joint work with Yijia Wang (Pitt)*

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# 1. Motivation via example applications

# Fast and slow states

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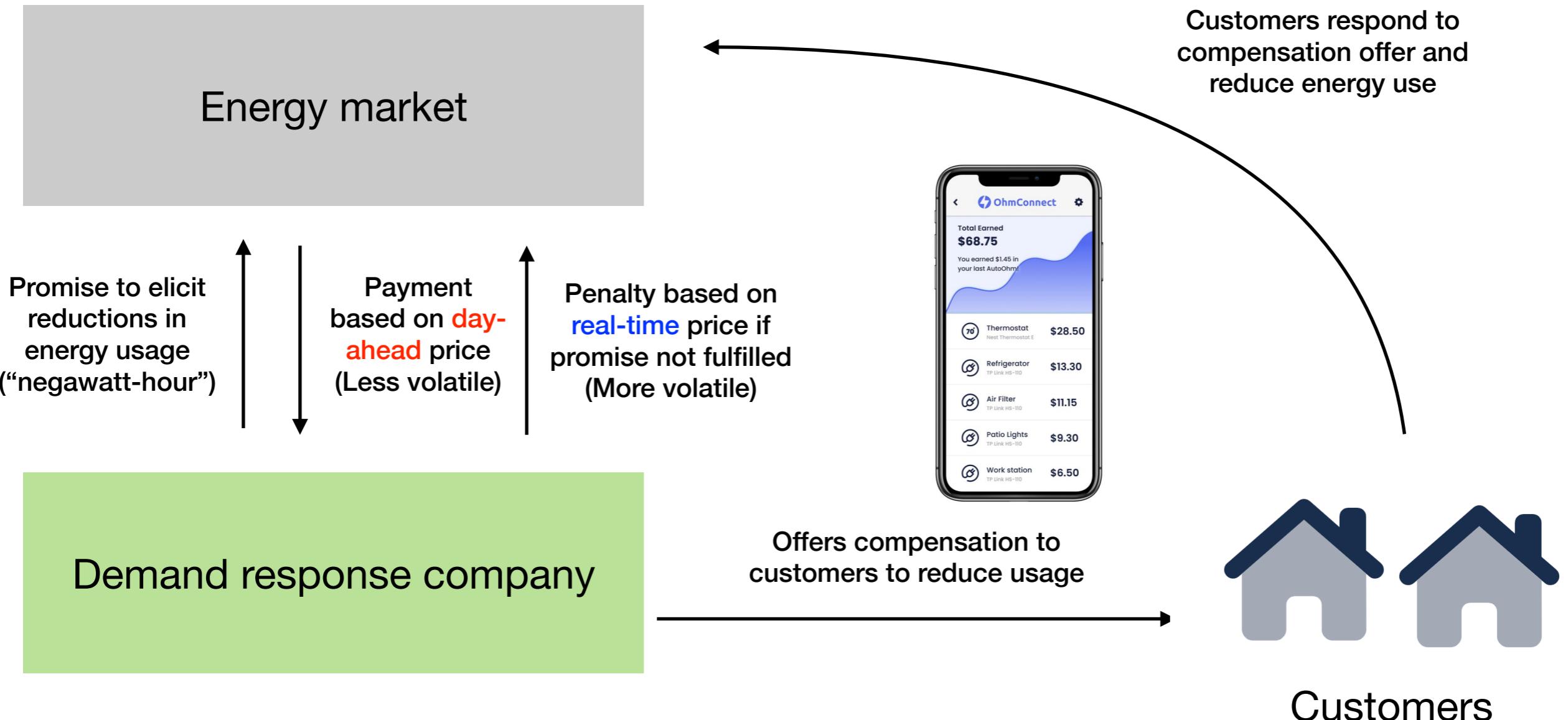


# Recommendation systems

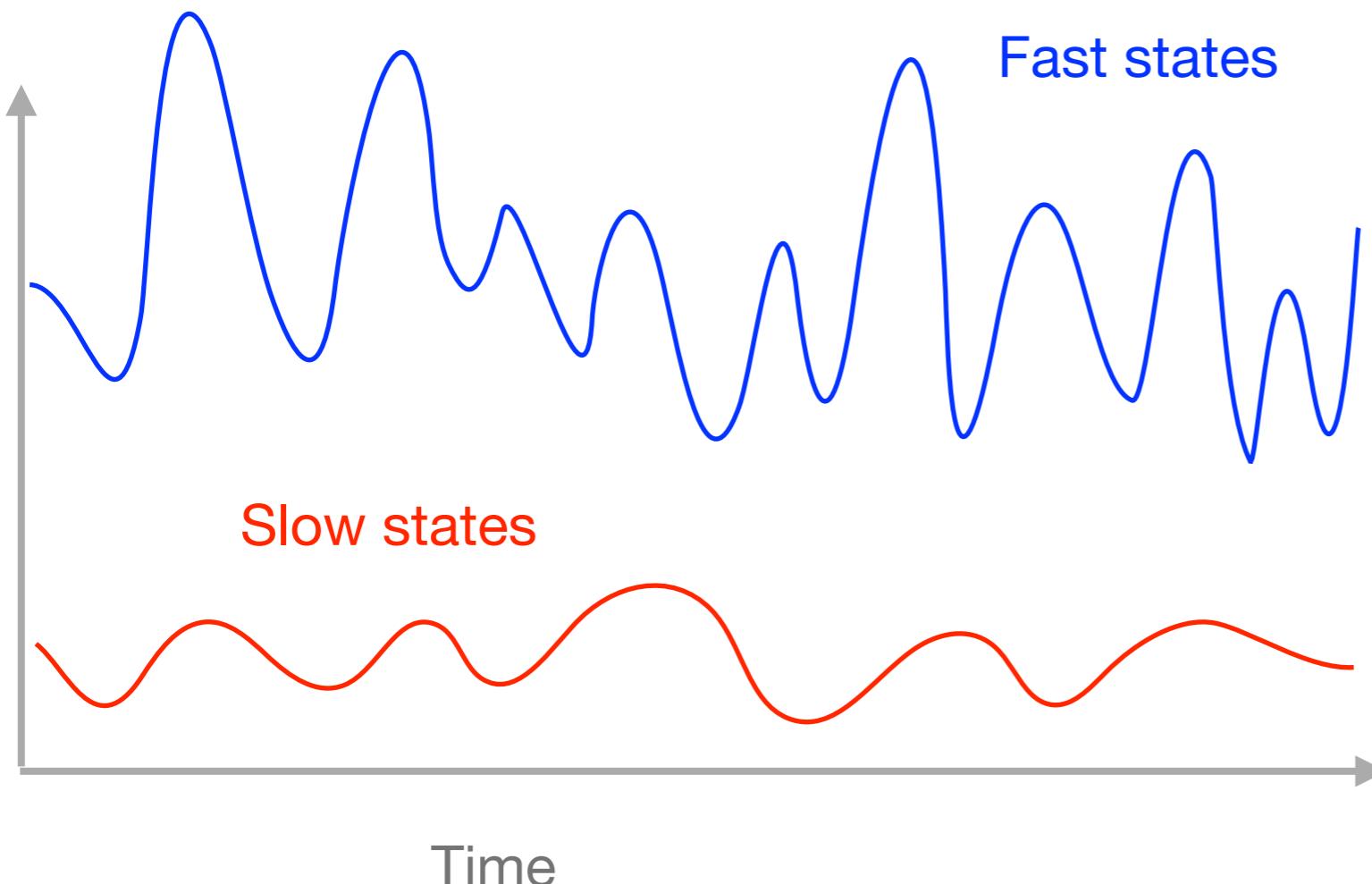
- Consider a recommendation system setting where:
  - Users return (i.e., log on) to the platform more often if they are seeing **content** that is *interesting* (Sumida & Zhou, 2023)
  - Users return to the platform more often if they have a *diverse recent content history*.
  - Users **interests** can shift over time as a function of the content they see.



# Demand response (shifting electricity demand)



# What do they have in common?



**Fast states** from examples:

- Real-time prices
- Recent content history

*Shorter timescales*

**Slow states** from examples:

- Day-ahead prices
- Underlying user interests

*Longer timescales*

# Current practice when modeling a new problem

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- While modeling an MDP, additional state variables is expensive:
  - Each iteration of value iteration  $\mathcal{O}(S^2A)$
- **What do practitioners do (anecdotally)?**
  - If a state is deemed a “slow state” (contexts, environmental variables, etc), they might be *ignored/omitted*
    - e.g. assume costs are deterministic, demand is stationary, weather doesn’t change
- **This work:** A *compromise* between computational tractability and fully ignoring the slow state
  - We propose to *periodically* ignore slow states
  - We give evidence and argue that completely omitting slow states from the decision model is often not a viable heuristic



## 2. Fast-slow Markov decision processes

# Fast-slow Markov decision processes

- A  $\gamma$ -discounted, infinite horizon MDP:
  - States  $s \in \mathcal{S}$
  - Actions  $a \in \mathcal{A}$
  - Rewards  $r(s, a) \in [0, r_{\max}]$
  - Transition function
    - $s_{t+1} = f(s_t, a_t, w_{t+1}), w_{t+1} \in \mathcal{W}$
- Fast-slow MDP:
  - States  $s = (x, y) \in \mathcal{S} = (\mathcal{X} \times \mathcal{Y})$
  - Actions  $a \in \mathcal{A}$
  - Rewards  $r(s, a) \in [0, r_{\max}]$
  - Transition function
    - $x_{t+1} = f_{\mathcal{X}}(s_t, a_t, w_{t+1})$
    - $y_{t+1} = f_{\mathcal{Y}}(s_t, a_t, w_{t+1})$

Main assumption (“fast-slow property”):

$$\|y - f_{\mathcal{Y}}(x, y, a, w)\|_2 \leq d_{\mathcal{Y}} \quad \text{and} \quad \|x - f_{\mathcal{X}}(x, y, a, w)\|_2 \leq \alpha d_{\mathcal{Y}}.$$

Lipschitz assumptions (let  $U^*(s)$  be the optimal value function):

$$|r(s, a) - r(s', a')| \leq L_r \|(s, a) - (s', a')\|_2,$$

$$\|f(s, a, w) - f(s', a', w)\|_2 \leq L_f \|(s, a) - (s', a')\|_2,$$

$$\|U^*(s) - U^*(s')\|_2 \leq L_U \|s - s'\|_2.$$

### 3. Hierarchical reformulation

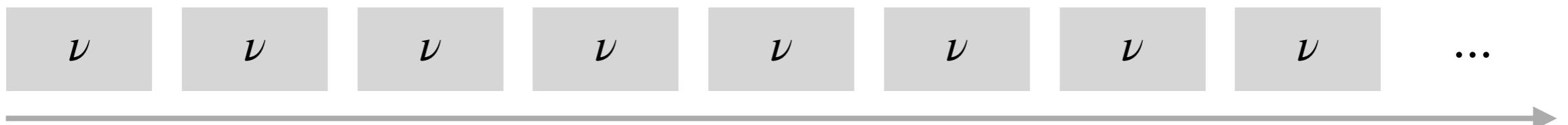
# Hierarchical reformulation (of any MDP)

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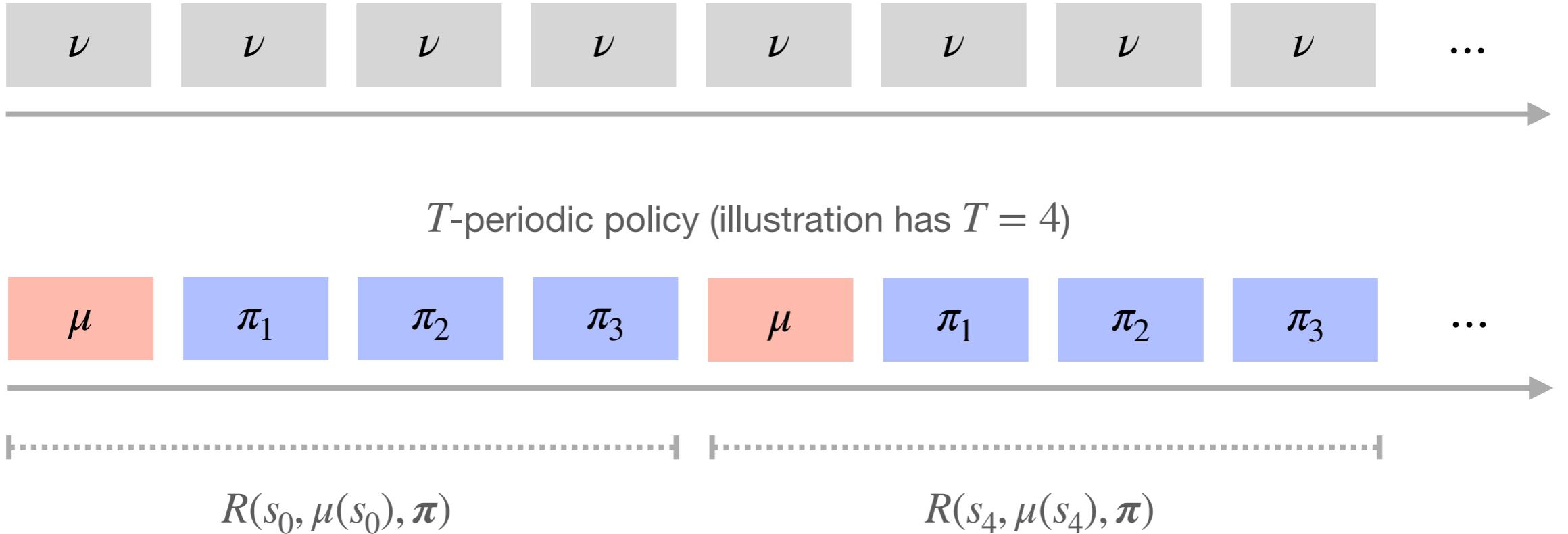
- A hierarchical reformulation is at the basis of our proposed approach
- Consider an MDP  $\langle \mathcal{S}, \mathcal{A}, \mathcal{W}, f, r, \gamma \rangle$
- Let  $\nu : \mathcal{S} \rightarrow \mathcal{A}$  be a stationary policy
- The value function is

$$U^\nu(s) = \mathbb{E} \left[ \sum_{t=0}^{\infty} \gamma^t r(s_t, \nu) \mid s_0 = s \right] = r(s, \nu) + \gamma \mathbb{E}[U^\nu(s')]$$

- The policy can be thought of as  $(\nu, \nu, \dots)$ :



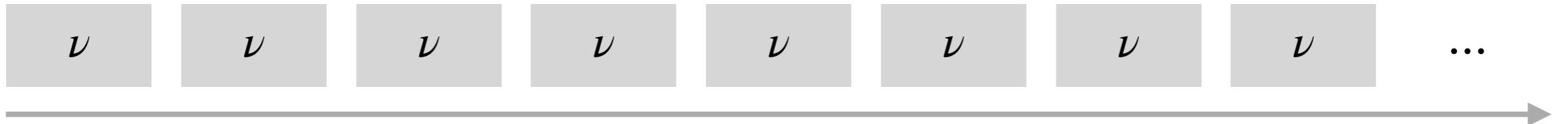
# Hierarchical reformulation (of any MDP)



- Given a  $T$ -periodic policy  $(\mu, \pi) = (\mu, \pi_1, \dots, \pi_{T-1})$ ,  $T$ -horizon reward is

$$R(s_0, \mu(s_0), \pi) = r(s_0, \mu) + \sum_{t=1}^{T-1} \gamma^t r(s_t, \pi_t)$$

# Hierarchical reformulation (of any MDP)



$T$ -periodic policy (illustration has  $T = 4$ )



- Bellman equations of the base model and its hierarchical reformulation are:

$$U^*(s_0) = \max_a r(s, a) + \gamma \mathbb{E}[U^*(s_1)]$$
$$\bar{U}^*(s_0) = \max_{(\mu, \pi)} \mathbb{E}[R(s_0, \mu(s_0), \pi) + \gamma^T \bar{U}^*(s_T)]$$

How can we take advantage of this?

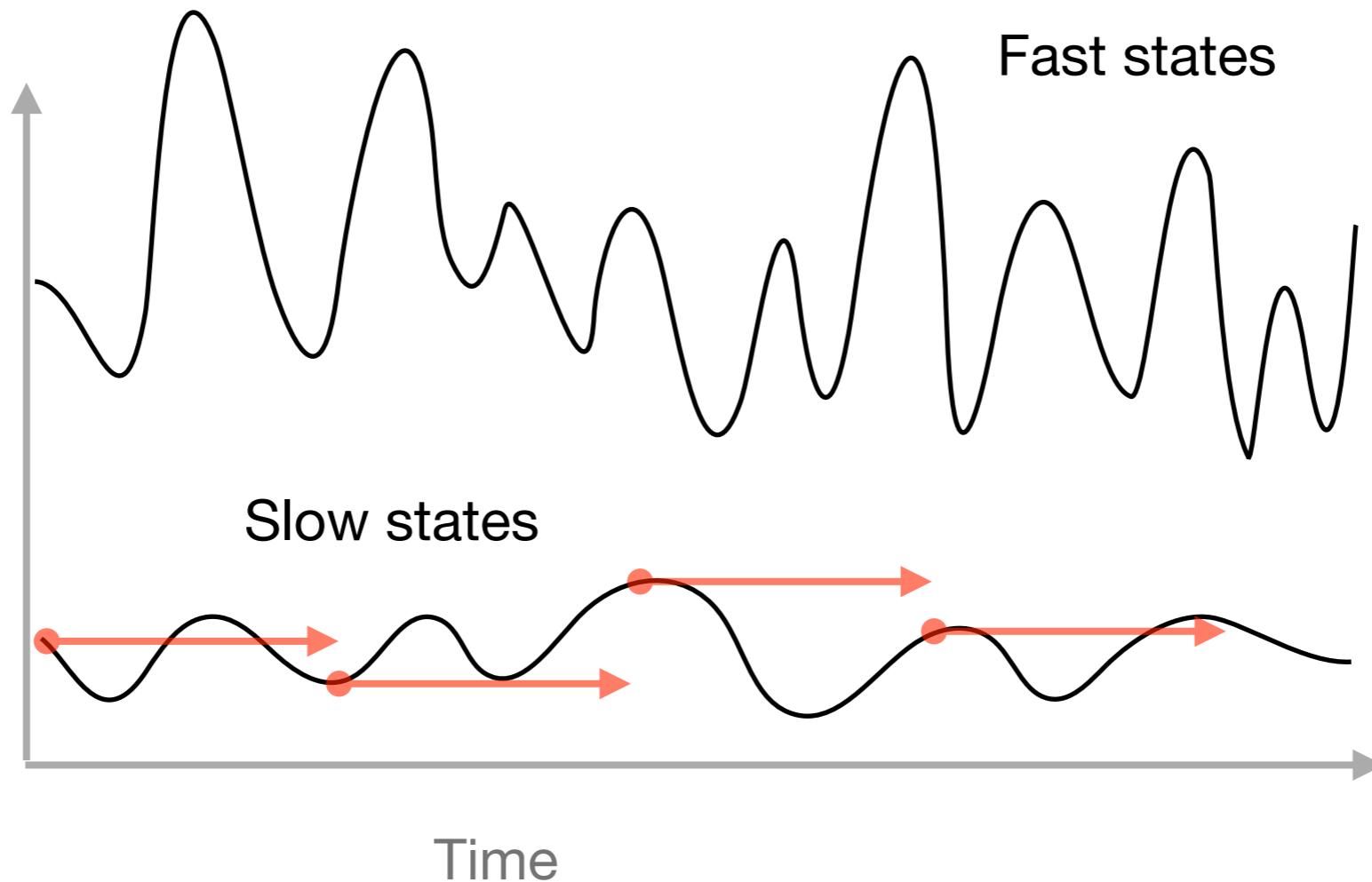
**Proposition.** The optimal values are equal:  $U^*(s) = \bar{U}^*(s)$ . Therefore, we can use the hierarchical reformulation as a basis for our approximation.

## 4. Frozen-state approximation and its regret

# Frozen-state approximation



**Main idea:** Since slow states don't change much, let's **freeze** them for some number ( $T$ ) of periods. Easier sub-problem with a smaller state space.



# Frozen-state approximation



**Main idea:** Since slow states don't change much, let's **freeze** them for some number ( $T$ ) of periods. Easier sub-problem with a smaller state space.

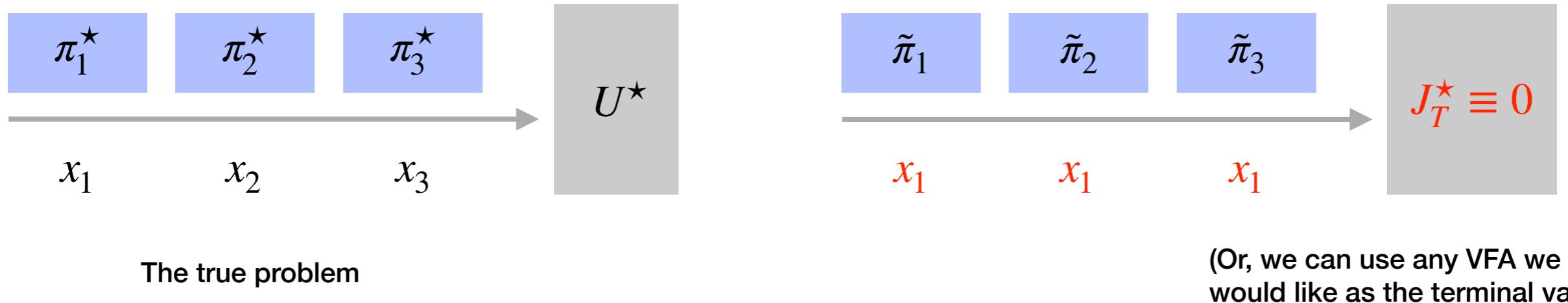
## Implementation

1. At  $t = 0$ , take a “upper-level” action (using  $\tilde{\mu}$ ), i.e., an action that considers the  $\gamma^T$  timescale
2. At  $t = 1$ , observe slow state and pretend it is frozen until  $t = T$  and that  $t = T$  is the end of the horizon
3. Solve this *easier* lower-level finite horizon problem.
4. Execute this  $T$ -period lower-level policy  $(\tilde{\pi}_1, \tilde{\pi}_2, \dots, \tilde{\pi}_{T-1})$  in the real system
5. Repeat

## Computation

- **Pre-compute** finite-horizon lower-level policy with frozen slow states
- Re-use pre-computed lower-level policy to solve infinite-horizon upper-level problem, which **takes advantage of**  $\gamma^T$

# Frozen-state, lower-level problem



## Frozen-state lower-level MDP

$$J_1^*(x, y) = \max_{\tilde{\pi}} \mathbb{E} \left[ \sum_{t=1}^{T-1} \gamma^{t-1} r(\textcolor{red}{x}_1, y_t, \tilde{\pi}_t) \mid (x_1, y_1) = (x, y) \right]$$

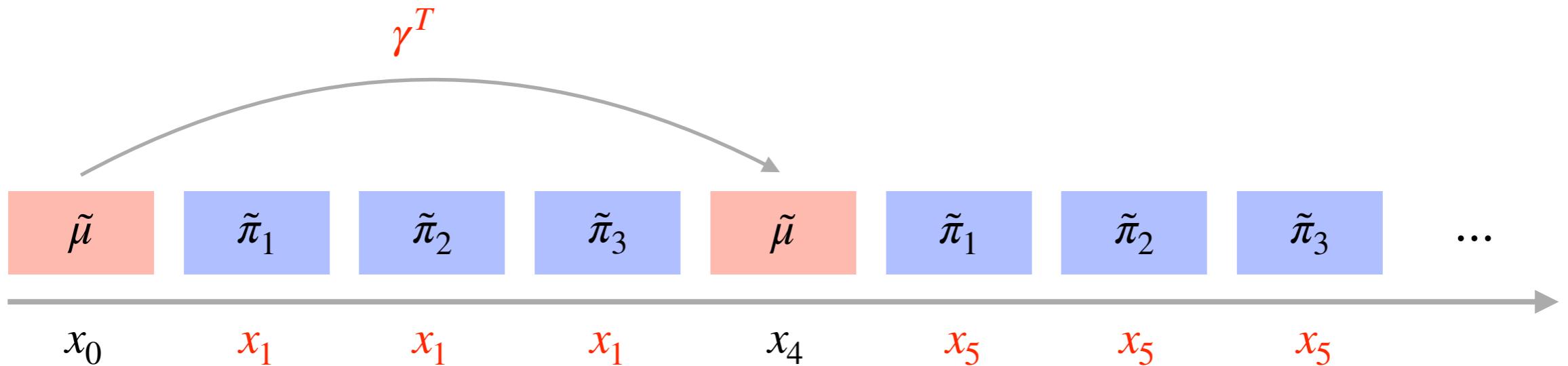
$$J_t^*(x, y) = \max_a r(x, y, a) + \gamma \mathbb{E}[J_{t+1}^*(x, y')], \quad \textcolor{red}{J}_T^* \equiv 0$$

$$\tilde{\pi}_t^*(x, y) = \operatorname{argmax}_a r(x, y, a) + \gamma \mathbb{E}[J_{t+1}^*(x, y')].$$

## Computational benefits

- Small number of successor states (since slow state is frozen)
  - $\mathcal{O}(S^2 A) \rightarrow \mathcal{O}(XY^2 A)$
- Independent across  $x$
- Independent from upper-level problem (replaced  $U^*$  by 0)

# Frozen-state, upper-level problem



## Frozen-state upper-level MDP

Let  $(\tilde{\pi}^{\star}, J_1^{\star})$  be the optimal policy/value of the lower-level problem.

$$\tilde{R}(s_0, a, J_1^{\star}) = r(s_0, a) + \gamma J_1^{\star}(f(s_0, a, w))$$

$$V^{\star}(s_0, J_1^{\star}, \tilde{\pi}^{\star}) = \max_a \mathbb{E}[\tilde{R}(s_0, a, J_1^{\star}) + \gamma^T V^{\star}(s_T, J_1^{\star}, \tilde{\pi}^{\star})] \text{ (transitions according to } \tilde{\pi}^{\star})$$

After solving both levels, let  $(\tilde{\mu}^{\star}, \tilde{\pi}^{\star})$  be the solution of the frozen-state approximation.

In the exact reformulation, we were maximizing over policies, now it is just a single action.

# Per-cycle reward approximation error

**Proposition.** The difference between true and approximate  $T$ -horizon rewards:

$$\begin{aligned}
 & \left| \underbrace{\mathbb{E}[R(s_0, a, \pi^*)]}_{\text{True}} - \underbrace{\mathbb{E}[\tilde{R}(s_0, a, J_1^*)]}_{\text{Frozen}} \right| \\
 & \leq \underbrace{\alpha d_{\mathcal{Y}} \left( L_r \sum_{i=1}^{T-2} \gamma^i \sum_{j=0}^{i-1} L_f^j \right)}_{\text{error from freezing}} + \underbrace{\gamma^{T-1} L_U \left[ \alpha d_{\mathcal{Y}} \sum_{j=0}^{T-2} L_f^j + \gamma d_{\mathcal{Y}} (\alpha + 2)(T - 1) \right]}_{\text{end of horizon error}}
 \end{aligned}$$

**Main ideas.**

1.  $\mathbb{E}[R(x_0, y_0, a, \pi^*)] = \mathbb{E}\left[r(x_0, y_0, a) + \gamma (H^{T-1} U^*)(x_1, y_1) - \gamma^T U^*(x_T, y_T)\right]$

where  $(HU)(x, y) = \max_a r(x, y, a) + \gamma \mathbb{E}[f(s, a, w))]$  (true Bellman operator)

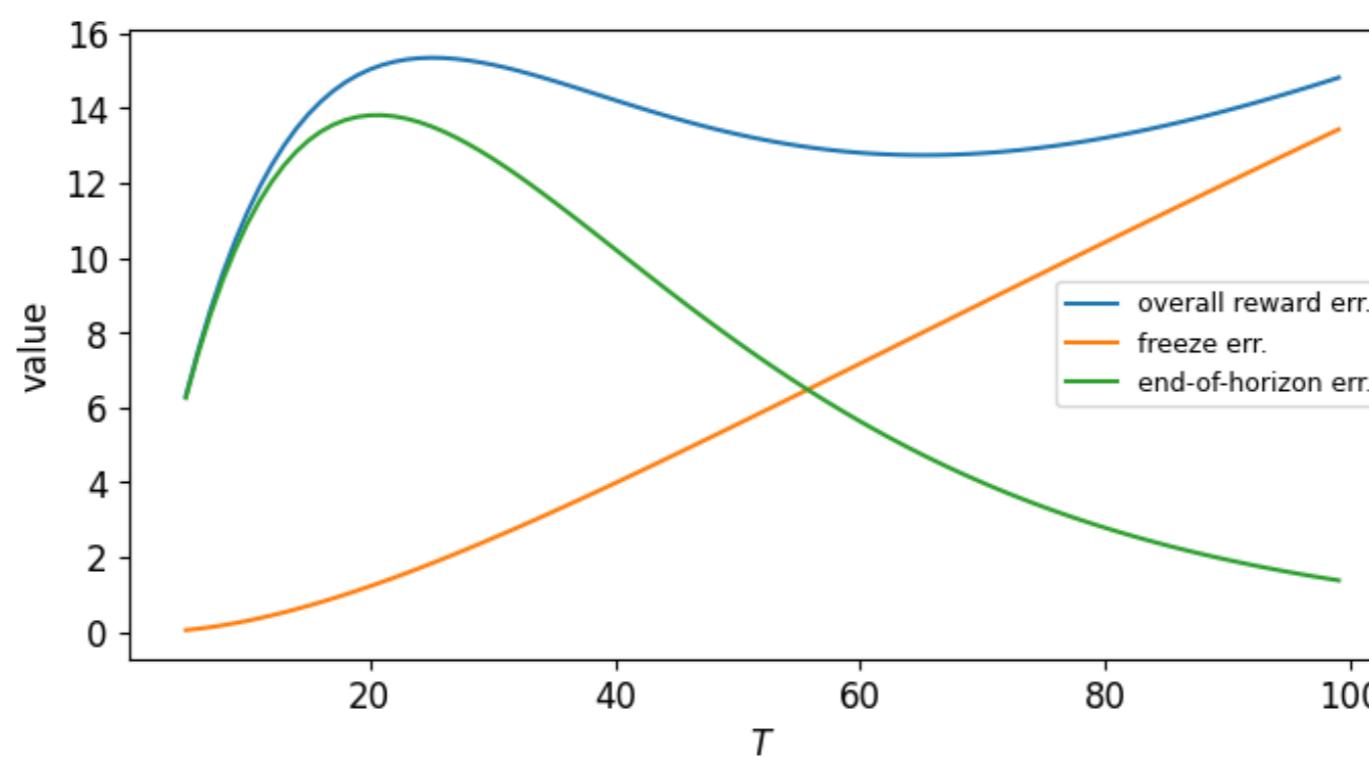
2.  $\mathbb{E}[\tilde{R}(x_0, y_0, a, J_1^*)] = r(x_0, y_0, a) + \gamma (\tilde{H}^{T-1} \mathbf{0})(x_1, y_1)$

where  $(\tilde{H}J_{t+1})(x, y) = \max_a r(x, y, a) + \gamma \mathbb{E}[J_{t+1}(x, f_{\mathcal{Y}}(x, y, a, w))]$  (frozen Bellman operator)

# Per-cycle reward approximation error

**Proposition.** The difference between true and approximate  $T$ -horizon rewards:

$$\left| \underbrace{\mathbb{E}[R(s_0, a, \pi^\star)] - \mathbb{E}[\tilde{R}(s_0, a, J_1^\star)]}_{\text{True}} \right| \leq \underbrace{\alpha d_{\mathcal{Y}} \left( L_r \sum_{i=1}^{T-2} \gamma^i \sum_{j=0}^{i-1} L_f^j \right)}_{\text{error from freezing}} + \underbrace{\gamma^{T-1} L_U \left[ \alpha d_{\mathcal{Y}} \sum_{j=0}^{T-2} L_f^j + \gamma d_{\mathcal{Y}} (\alpha + 2)(T - 1) \right]}_{\text{end of horizon error}}$$



## 5. A new algorithm: *Frozen-state value iteration*

# Standard value iteration on the base model

**Recall:** Given an MDP and Bellman operator  $H$ , where  $(HU)(s) = \max_a r(s, a) + \gamma \mathbb{E}U(f(s, a, w))$ , the *value iteration* algorithm is  $U^k = H^k U^0$

. Convergence to optimal value function:  $\lim_{t \rightarrow \infty} H^t U = U^*$  for any initial estimate  $V$

.  $\|U^{\nu_k} - U^*\|_\infty \leq \frac{2r_{\max}\gamma^{k+1}}{(1-\gamma)^2}$ , where  $\nu^k(s) = \operatorname{argmax}_a r(s, a) + \gamma \mathbb{E}[U^k(f(s, a, w))]$



Depends on

.  $\|U^k - U^*\|_\infty \leq \gamma^k \|U^0 - U^*\|_\infty$

.  $\|U^0 - U^*\|_\infty \leq \frac{r_{\max}}{1-\gamma}$

.  $\|U^{\nu_k} - U^*\|_\infty \leq \frac{2\|U^k - U^*\|_\infty}{1-\gamma}$

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## Algorithm 1: Exact VI for the Base Model

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**Input:** Initial values  $U_0$ , number of iterations  $k$ .

**Output:** Approximation to the optimal policy  $\nu^k$ .

1 **for**  $i = 1, 2, \dots, k$  **do**

2   **for**  $s$  in the state space  $\mathcal{S}$  **do**

3      $U^i(s) = \max_a r(s, a) + \gamma \mathbb{E}[U^{i-1}(f(s, a, w))]$ .

4   **end**

5 **end**

6 **for**  $s$  in the state space  $\mathcal{S}$  **do**

7      $\nu^k(s) = \operatorname{argmax}_a r(s, a) + \gamma \mathbb{E}[U^k(f(s, a, w))]$ .

8 **end**

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# Frozen-state value iteration (FSVI)

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**Algorithm 2:** Frozen-State Value Iteration (FSVI)

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**Input:** Initial values  $J_T^* \equiv 0$  and  $V^0$ , number of iterations  $k$ .

**Output:** Approximation of the  $T$ -periodic frozen-state policy  $(\tilde{\mu}^k, \tilde{\pi}^*)$  and  $J_1^*$ .

```

1 for  $t = T - 1, T - 2, \dots, 1$  do
2   for each slow state  $x \in \mathcal{X}$  do
3     for each fast state  $y \in \mathcal{Y}$  do
4        $J_t^*(x, y) = \max_a r(x, y, a) + \gamma \mathbb{E}[J_{t+1}^*(x, f_y(x, y, a, w))]$ .
5        $\tilde{\pi}_t^*(x, y) = \arg \max_a r(x, y, a) + \gamma \mathbb{E}[J_{t+1}^*(x, f_y(x, y, a, w))]$ .
6     end
7   end
8 end
9
9 for  $i = 1, 2, \dots, k$  do
10  for  $s_0 = (x_0, y_0)$  in the state space  $\mathcal{X} \times \mathcal{Y}$  do
11     $V^i(x_0, y_0, J_1^*, \tilde{\pi}^*) = \max_a \mathbb{E}[\tilde{R}(s_0, a, J_1^*) + \gamma^T V^{i-1}(x_T, y_T, J_1^*, \tilde{\pi}^*)]$ .
12  end
13 end
14
14 for  $s_0 = (x_0, y_0)$  in the state space  $\mathcal{X} \times \mathcal{Y}$  do
15   $\tilde{\mu}^k(x_0, y_0) = \arg \max_a \mathbb{E}[\tilde{R}(s_0, a, J_1^*) + \gamma^T V^k(x_T, y_T, J_1^*, \tilde{\pi}^*)]$ .
16 end

```

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**Solving the lower level incurs a one time fixed cost**

Pre-compute lower-level problem, a finite-horizon DP:

- To solve lower-level DP:  $\mathcal{O}(XY^2AT)$
- To compute multi-step transition:  $\mathcal{O}(S^2T)$

Upper-level problem (infinite-horizon VI on slow-timescale MDP with  $\gamma^T$  discounting):

- Per upper-level VI iteration:  $\mathcal{O}(S^2A)$

$\mathcal{O}(S^2A)$  per iteration is the same as Base VI, but now the discount factor is  $\gamma^T$  instead of  $\gamma$ !

# Regret of a periodic policy $(\mu, \pi)$

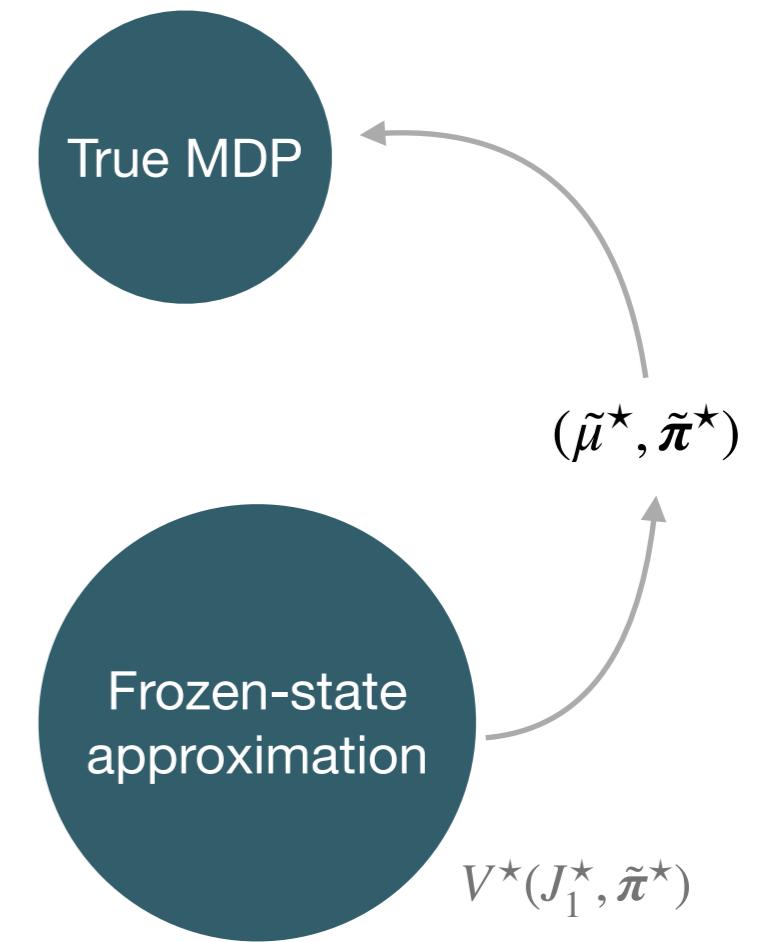
**Definition.** Suppose the optimal policy is  $\nu^*$ . The regret is

$$\mathcal{R}(s, \mu, \pi) = U^{\nu^*}(s) - \bar{U}^{\mu}(s, \pi) = \bar{U}^*(s) - \bar{U}^{\mu}(s, \pi) \quad \text{and} \quad \mathcal{R}(\mu, \pi) = \max_s \mathcal{R}(s, \mu, \pi),$$

where we have used the equivalence between the base and hierarchical formulations.

## Remarks:

- We always measure regret with respect to the *true MDP*.
  - Although  $(\mu, \pi)$  is computed *with the help of frozen states*, it is evaluated in the original MDP with true dynamics.
- Consider  $\mathcal{R}(\tilde{\mu}^*, \tilde{\pi}^*)$ , notice that  $V^*(J_1^*, \tilde{\pi}^*)$  does not directly enter the regret definition.
  - It is the optimal value of the approximation, but doesn't reflect the performance of  $(\tilde{\mu}^*, \tilde{\pi}^*)$  in the true model.



# Main idea behind regret analysis

## Lemma (Approximation to FSVI).

- Suppose we *approximately* solve the lower-level problem and obtain  $\pi, J_1$ , instead of the optimal solutions  $\pi^*, U^*$ .
- Suppose we approximately solve the upper-level problem and obtain  $V$  instead of  $V^*(J_1, \pi)$ .
- Let  $\mu$  be greedy with respect to both  $J_1$  and  $V$ :
  - $\mu(s_0) = \operatorname{argmax}_{a \in \mathcal{A}} \mathbb{E}[\tilde{R}(s_0, a, J_1) + \gamma^T V(s_T(a, \pi))]$ .
- Then,

$$\mathcal{R}(\mu, \pi) \leq \left( \frac{2\gamma^T}{(1 - \gamma^T)^2} + \frac{2}{1 - \gamma^T} \right) \epsilon_r(\pi^*, J_1) + \left( \frac{2\gamma^{2T}}{(1 - \gamma^T)^2} + \frac{2\gamma^T}{1 - \gamma^T} \right) L_U d(a, d_y, T) + \frac{2\gamma^T}{1 - \gamma^T} \|V^*(J_1, \pi) - V\|_\infty.$$

End of horizon error

V-approximation error

Reward error

# Regret of FSVI

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**Theorem.** The regret of FSVI after  $k$  upper-level iterations is:

$$\begin{aligned}\mathcal{R}(\mu, \boldsymbol{\pi}) \leq & \left( \frac{2\gamma^T}{(1 - \gamma^T)^2} + \frac{2}{1 - \gamma^T} \right) \epsilon_r(\boldsymbol{\pi}^\star, J_1) \\ & + \left( \frac{2\gamma^{2T}}{(1 - \gamma^T)^2} + \frac{2\gamma^T}{1 - \gamma^T} \right) L_U d(\alpha, d_{\mathcal{Y}}, T) + \frac{2r_{\max}\gamma^{(k+1)T}}{(1 - \gamma)(1 - \gamma^T)},\end{aligned}$$

which replaces the V-approximation error term with the VI error.

## Main question to answer using the regret analysis

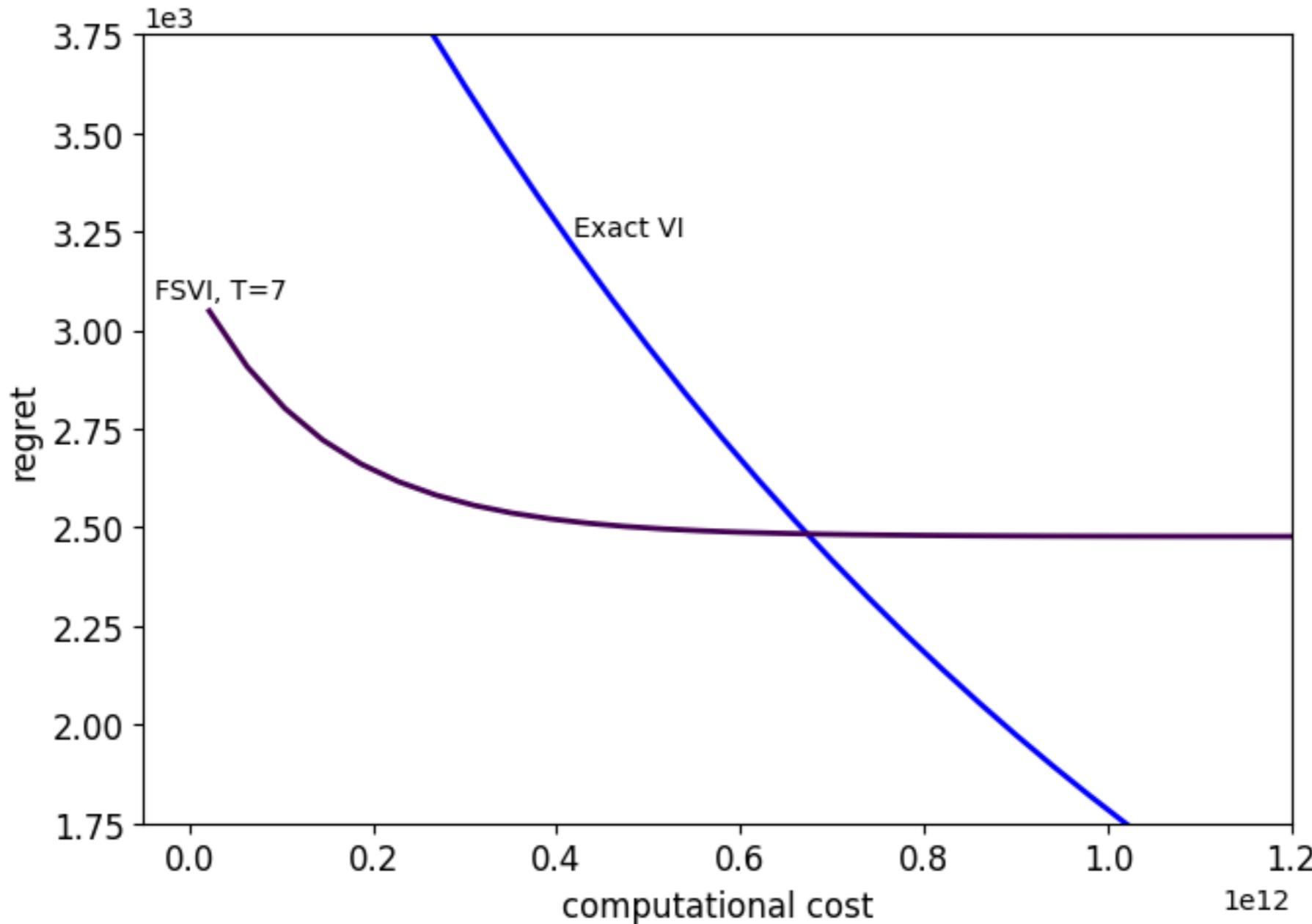
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*Given some computational budget, should we use the FSVI algorithm at all?*

*If so, how should we choose  $T$  (the number of periods to freeze) and  $k$  (number of upper-level steps iterations)?*

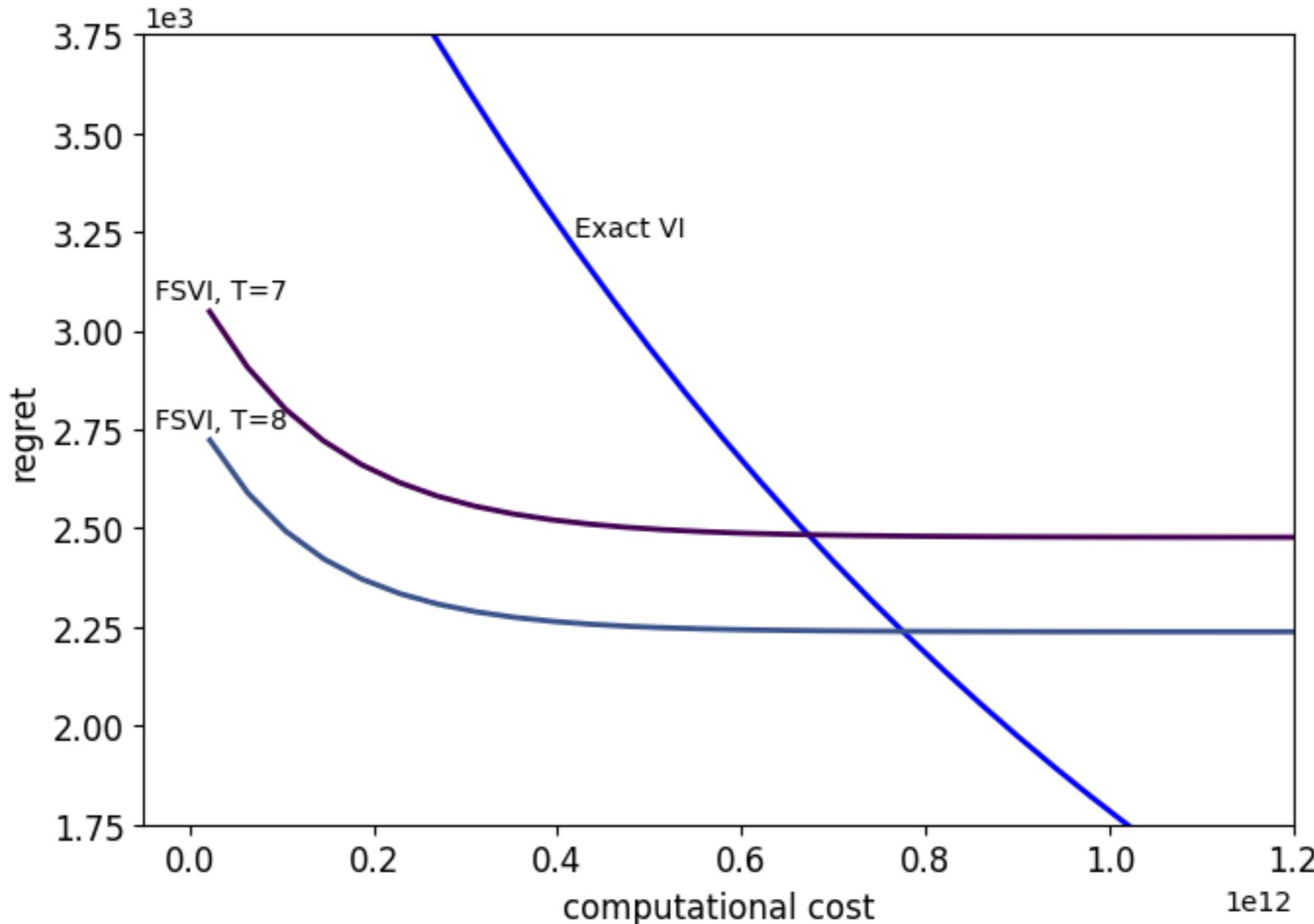
## How can we use the bound to choose T?

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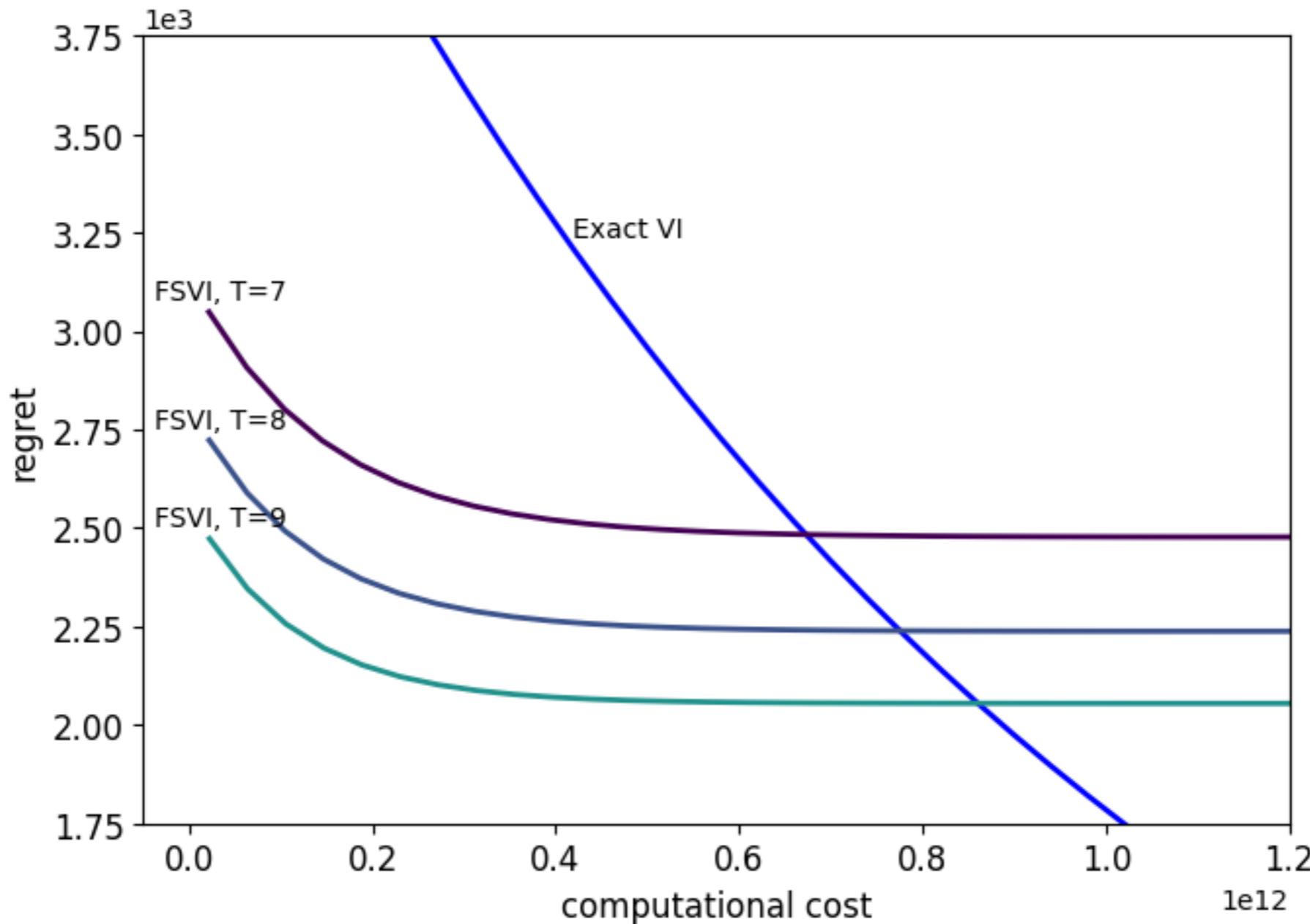
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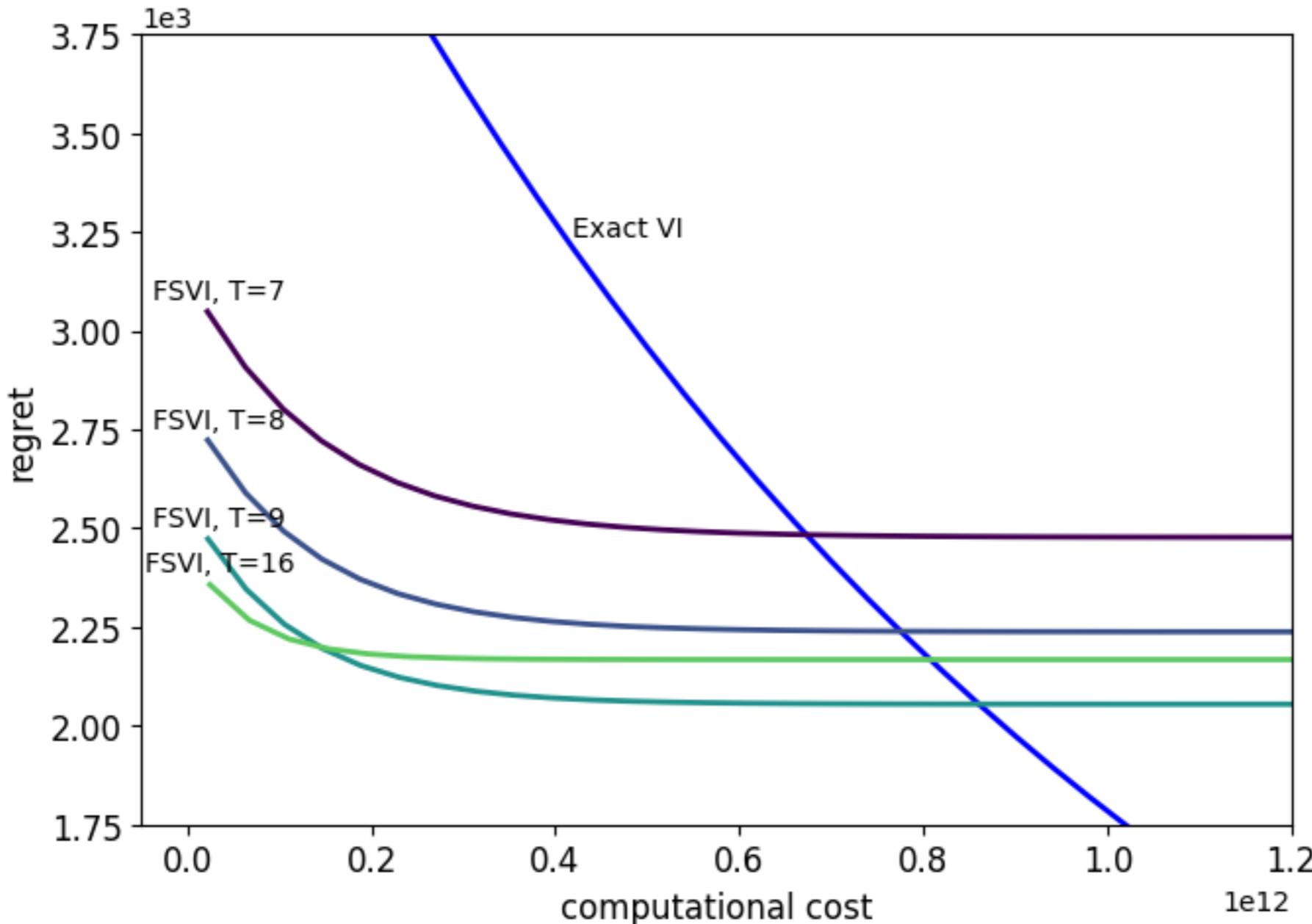
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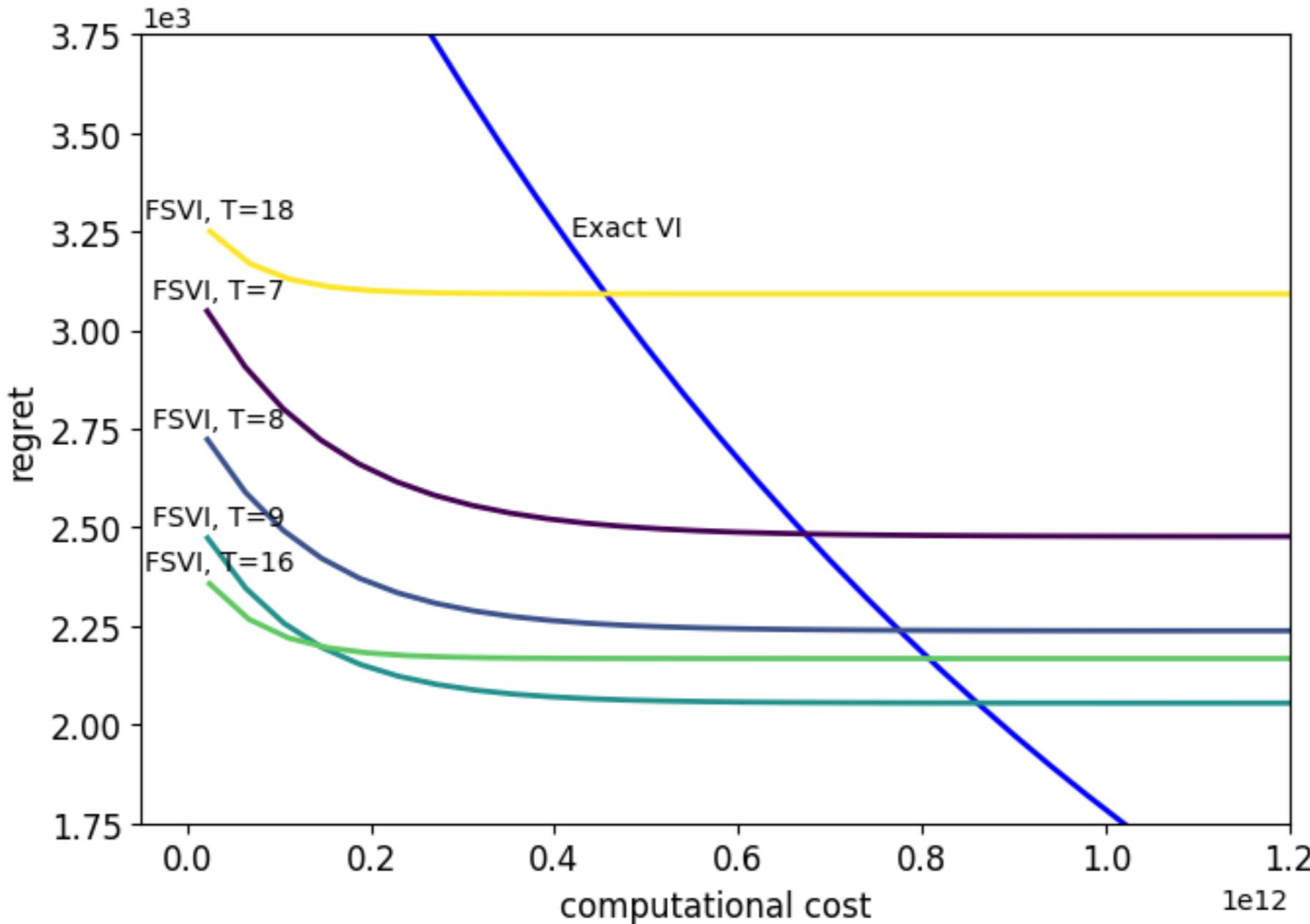
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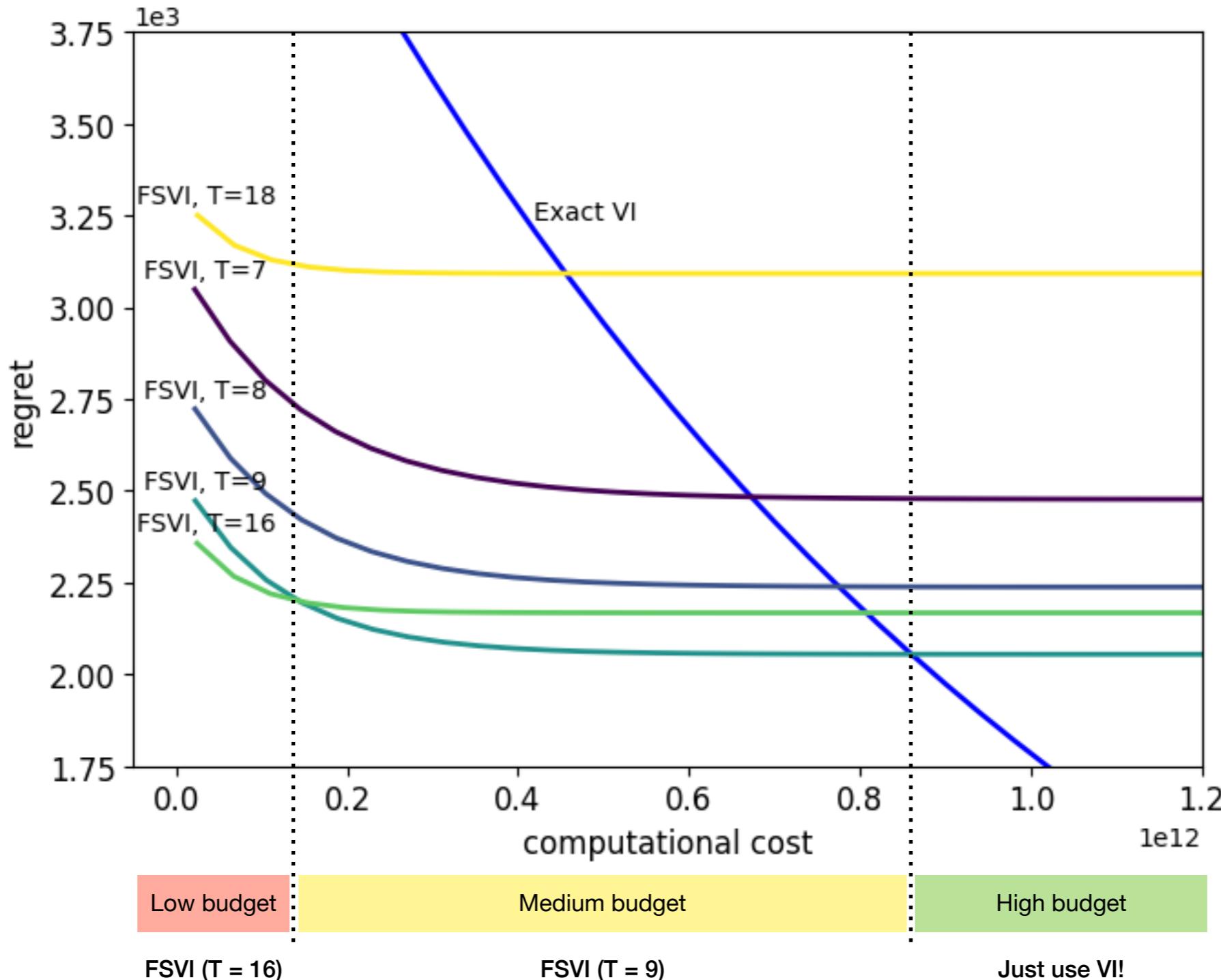


## How can we use the bound to choose T?

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# How can we use the bound to choose T?

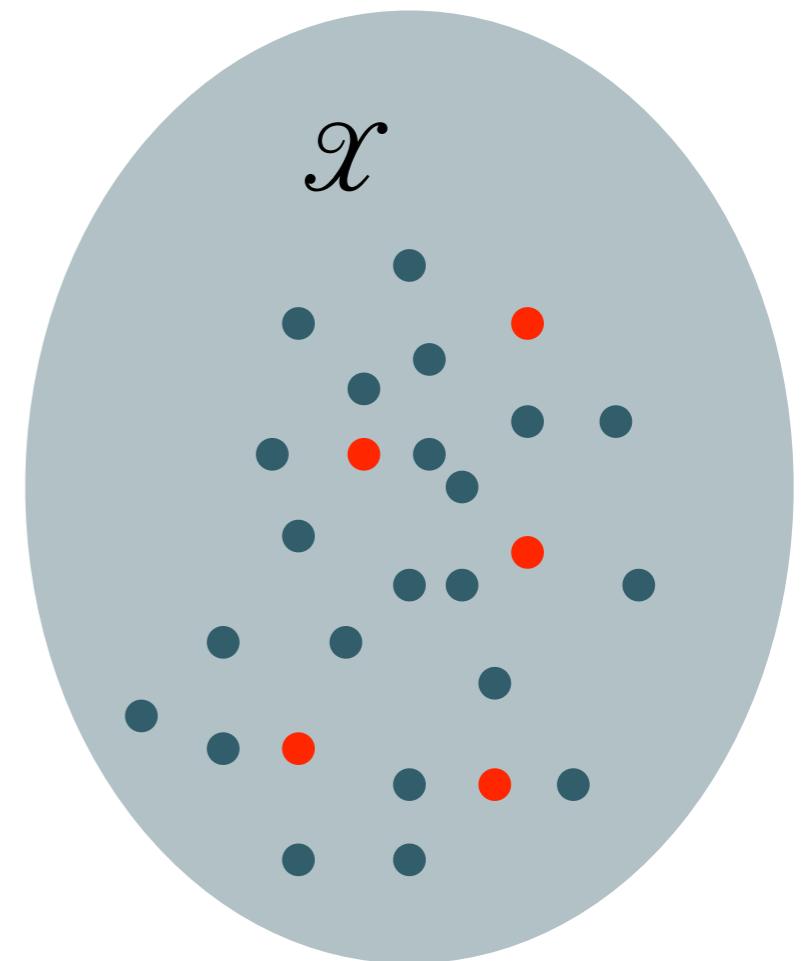


## 6. Extensions

## Extension: Nominal state version of FSVI

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- In FSVI, we have to solve the lower-level MDP for each  $x$
- We can further do approximations by solving the lower-level MDP for a few nominal states only
  - $\mathcal{O}(S^2A) \rightarrow \mathcal{O}(XY^2A) \rightarrow \mathcal{O}(X_{\text{nom}}Y^2A)$
- Later, extrapolate to nearby nominal states
- Theoretical results can be adapted given an additional assumption on the MDP rewards



# Extension: Scaling to larger state spaces using feature-based approximate value iteration

## Architecture:

- Consider  $M$  pre-selected states  $\tilde{\mathcal{S}} = \{s_1, s_2, \dots, s_M\}$ .
- Consider an  $M$ -dimensional feature vector  $\phi(s)$ , where  $\phi(s_m)$  are linearly independent.
- Assume there exists  $\gamma' \in [\gamma, 1)$  s.t. for any  $s$ , there exists  $\theta_m(s)$ , where
  - $\sum_m |\theta_m(s)| \leq 1$  and  $\phi(s) = \frac{\gamma'}{\gamma} \sum_{m=1}^M \theta_m(s) \phi(s_m)$ .
- Lower level:  $\hat{J}(s, \omega_t) = \phi^\top(s) \omega_t$ .
- Upper level:  $\hat{V}(s, \beta^k) = \phi^\top(s) \beta^k$ .
- Update procedure:
  1. Compute Bellman update at pre-selected states only:  $y(s_m)$ .
  2. Compute next parameter vector ( $\omega_{t-1}$  or  $\beta^{k+1}$ ) such that the updated value function evaluated at the pre-selected states is equal: e.g.,  $\hat{J}(s_m, \omega_{t-1}) = y(s_m)$ .

Limited “expansion” after going to parameter space and back:

$$\|(\Phi\Phi^\dagger)(J) - (\Phi\Phi^\dagger)(J')\|_\infty \leq \kappa \|J - J'\|_\infty \quad (\kappa = \gamma'/\gamma)$$



# Extension: Scaling to larger state spaces using feature-based approximate value iteration

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**Algorithm 4:** Frozen-State Approximate Value Iteration (FSAVI)

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**Input:**  $\tilde{\mathcal{S}} = \{s_1, s_2, \dots, s_M\}$ ,  $\phi$ , initial weights  $\omega_T = \beta_0 = \mathbf{0}$ , number of iterations  $k$ .

**Output:** Approximation of the  $T$ -periodic frozen-state policy  $(\hat{\mu}_{(\beta^k, \omega^*)}, \hat{\pi}_{\omega^*})$  and  $\hat{J}_1(\omega^*)$

```
1 for  $t = T - 1, T - 2, \dots, 1$  do
2   for each pre-selected state  $s = (x, y) \in \tilde{\mathcal{S}}$  do
3      $J_t(x, y) = \max_a r(x, y, a) + \gamma \mathbb{E}[\hat{J}_{t+1}(x, f_{\mathcal{Y}}(x, y, a, w), \omega_{t+1})]$ .
4   end
5   Set remaining entries of  $J_t$  to zero. Update parameter vector:  $\omega_t^* = \Phi^\dagger J_t$ .
6 end
7 Let  $\hat{\pi}_{\omega^*}$  be greedy with respect to  $\hat{J}_t(\omega_t^*) = \Phi \omega_t^*$ , similar to (23).
8 for  $i = 1, 2, \dots, k$  do
9   for each pre-selected state  $s_0 \in \tilde{\mathcal{S}}$  do
10     $V^i(s_0) = \max_a \mathbb{E}[\tilde{R}(s, a, \hat{J}_1(\omega_1^*)) + \gamma^T \hat{V}(s_T(a, \tilde{\pi}_{\text{avi}}), \beta_{i-1})]$ .
11    Set remaining entries of  $V^i$  to zero. Update parameter vector:  $\beta_i = \Phi^\dagger V^i$ .
12  end
13 end
14 for  $s_0$  in the state space  $\mathcal{S}$  do
15    $\hat{\mu}_{(\beta^k, \omega^*)}(s_0) = \arg \max_a \mathbb{E}[\tilde{R}(s_0, a, \hat{J}_1(\omega_1^*)) + \gamma^T \hat{V}(s_T(a, \tilde{\pi}_{\omega^*}), \beta_k)]$ .
16 end
```

---

# Regret of FSAVI

**Theorem.** The regret of FSAVI after  $k$  upper-level iterations is:

$$\begin{aligned} \mathcal{R}(\mu, \boldsymbol{\pi}) \leq & \left( \frac{2\gamma^T}{(1-\gamma^T)^2} + \frac{2}{1-\gamma^T} \right) \epsilon_r(\boldsymbol{\pi}^*, \hat{J}_1(\boldsymbol{\omega}_1^*)) \\ & + \left( \frac{2\gamma^{2T}}{(1-\gamma^T)^2} + \frac{2\gamma^T}{1-\gamma^T} \right) L_U d(\alpha, d_{\mathcal{Y}}, T) + \underbrace{\left( \frac{1+\kappa}{1-\kappa\gamma^T} \right) \epsilon_{\text{up}}}_{\|V_{\boldsymbol{\omega}^*}^* - \hat{V}(\boldsymbol{\beta}^*)\|_{\infty}} + \underbrace{(\kappa\gamma^T)^k \left( \frac{\kappa^2 - \kappa^2(\kappa\gamma)^{T+1}}{(1-\kappa\gamma^T)(1-\kappa\gamma)} \right) r_{\max}}_{\|\hat{V}(\boldsymbol{\beta}^*) - \hat{V}(\boldsymbol{\beta}_k)\|_{\infty}}, \end{aligned}$$

$$\text{which } \epsilon_r(\boldsymbol{\pi}^*, \hat{J}_1(\boldsymbol{\omega}_1^*)) = \epsilon_r(\boldsymbol{\pi}^*, J_1^*) + \underbrace{\left( \frac{1+\kappa}{1-\kappa\gamma} - \frac{(\kappa\gamma)^T(1+\gamma)}{\gamma - \kappa\gamma^2} \right) \epsilon_{\text{low}}}_{\|J_t^* - \hat{J}_t(\boldsymbol{\omega}_t^*)\|_{\infty}}.$$

Upper-level feature approximation error

Lower-level feature approximation error

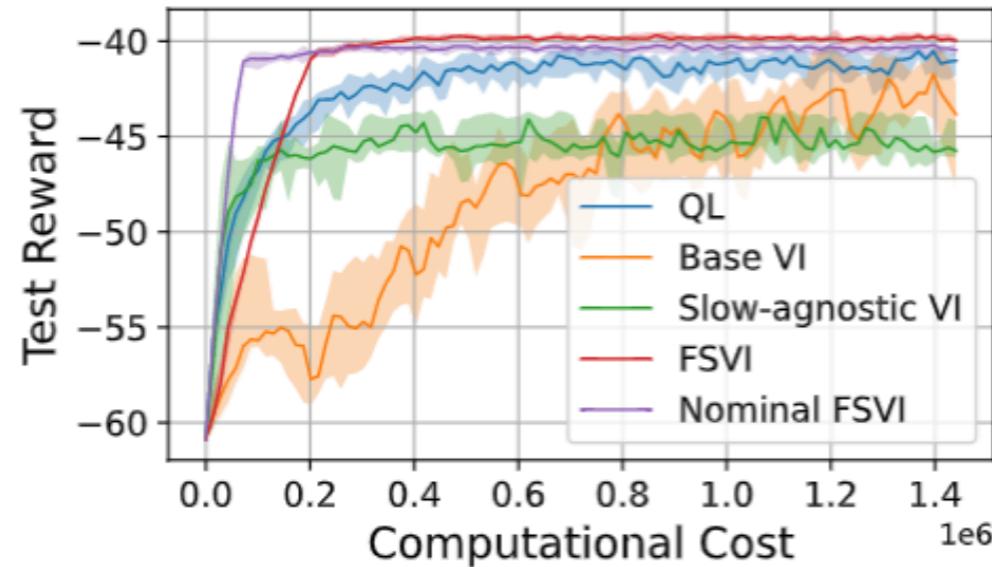
## 7. Numerical results

# Baseline algorithms

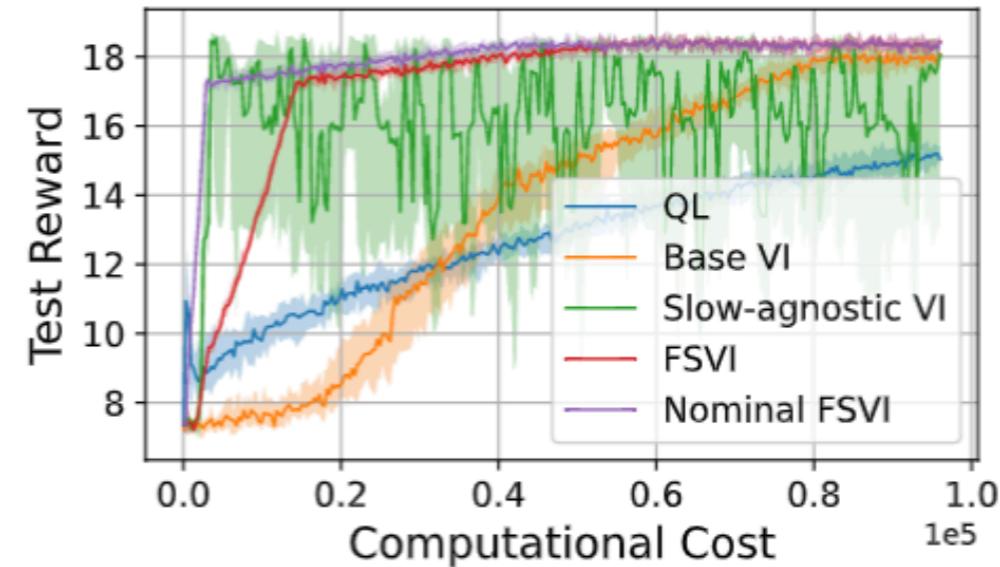
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- VI / AVI
- Slow-agnostic VI / AVI
  - Average over slow states during learning
  - Upon implementation, ignore slow state
- Q-learning (QL)
- Deep Q-networks (DQN)
- **Ours: FSVI / Nominal FSVI**
- **Ours: FSAVI / Nominal FSAVI**

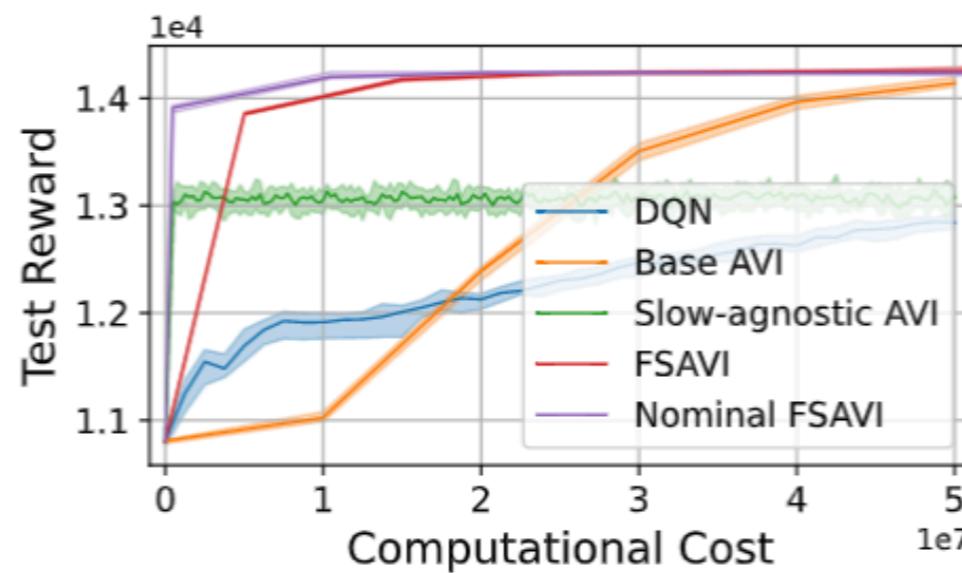
# Overall performance comparison



(a) Multi-class service allocation



(b) Restless two-armed bandit



(c) Energy demand response

# Questions

Please feel free to email me at [danielrjiang@gmail.com](mailto:danielrjiang@gmail.com) for additional comments and discussion.

