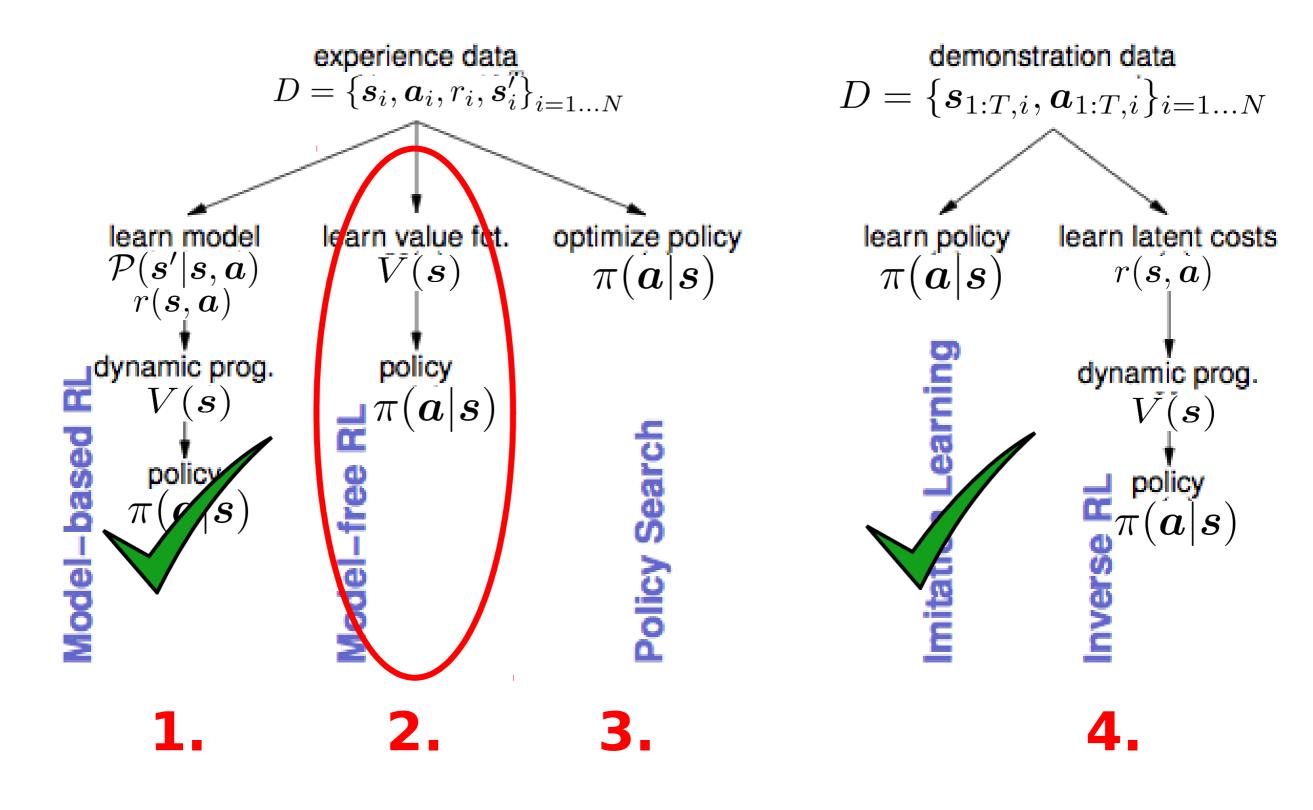
Reinforcement Learning Part 2: Value Function Methods

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The Bigger Picture: How to learn policies



CHRIS ATKESON Humanoids 2016

American election:

- Clinton was model-based, and used strong predictive models of who would vote and how they would vote.
- Trump did not use any models.



Often, learning a good model is too hard

- The optimization inherent in optimal control is prone to model errors, as the controller may achieve the objective only because model errors get exploited
- Optimal control methods based on linearization of the dynamics work only for moderately non-linear tasks
- Model-free approaches are needed that do not make any assumption on the structure of the model

Classical Reinforcement Learning:

Solve the optimal control problem by learning the value function, not the model!



1. Quick recap of dynamic programming

2. Reinforcement Learning with Temporal Differences

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Least-Squares Temporal Difference Learning

Fitted Q-Iteration

5. Robot Application: Robot Soccer

Final Remarks



Classical reinforcement learning is typically formulated for the infinite horizon objective

Infinite Horizon: maximize discounted accumulated reward

$$J_{\boldsymbol{\pi}} = \mathbb{E}_{\mu_0, \mathcal{P}, \boldsymbol{\pi}} \left[\sum_{t=0}^{\infty} \left[\gamma^t r(\boldsymbol{s}_t, \boldsymbol{a}_t) \right] \right]$$

$$\leq \gamma < 1$$

... discount factor

 $\left(\right)$

Value functions and State-Action Value Functions



Refresher: Value function and state-action value function can be computed iteratively

$$\begin{aligned} V^{\pi}(\boldsymbol{s}) &= \mathbb{E}_{\pi} \Big[r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \mathbb{E}_{\mathcal{P}} \left[V^{\pi}(\boldsymbol{s}') \right] \big| \boldsymbol{s} \Big] \\ &= \int \pi(\boldsymbol{a} | \boldsymbol{s}) \Big(r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \int \mathcal{P}(\boldsymbol{s}' | \boldsymbol{s}, \boldsymbol{a}) V^{\pi}(\boldsymbol{s}') d\boldsymbol{s}' \Big) d\boldsymbol{a} \end{aligned}$$

$$Q^{\pi}(\boldsymbol{s}, \boldsymbol{a}) = r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \mathbb{E}_{\mathcal{P}, \pi} \Big[Q^{\pi}(\boldsymbol{s}', \boldsymbol{a}') \big| \boldsymbol{s}, \boldsymbol{a} \Big]$$
$$= r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \int \mathcal{P}(\boldsymbol{s}' | \boldsymbol{s}, \boldsymbol{a}) \int \pi(\boldsymbol{a}' | \boldsymbol{s}') Q^{\pi}(\boldsymbol{s}', \boldsymbol{a}') d\boldsymbol{a}' d\boldsymbol{s}'$$

Bellman Equation of optimality

$$V^*(\boldsymbol{s}) = \max_{\boldsymbol{a}} \left(r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \mathbb{E}_{\mathcal{P}} \left[V^*(\boldsymbol{s}') \big| \boldsymbol{s}, \boldsymbol{a} \right] \right)$$

rating the Bellman Equation converges to the optimal value function V^* and is called **value iteration**

Alternatively we can also iterate Q-functions...

$$Q^*(\boldsymbol{s}, \boldsymbol{a}) = r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \mathbb{E}_{\mathcal{P}} \left[\max_{\boldsymbol{a}'} Q^*(\boldsymbol{s}', \boldsymbol{a}') \middle| \boldsymbol{s}, \boldsymbol{a} \right]$$



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Classical Reinforcement Learning

Updates the value function based on samples

 $\mathcal{D} = \{\boldsymbol{s}_i, \boldsymbol{a}_i, r_i, \boldsymbol{s}'_i\}_{i=1...N}$

We do not have a model and we do not want to learn it

Use the samples to update Q-function (or V-function)

Lets start simple:

Discrete states/actions 🗭 Tabular Q-function



Given a transition a_t, r_t, s_{t+1} , we want to update the V-function

 $V(s_t)$ Estimate of the current value:

$$\hat{V}(s_t) = r_t + \gamma V(s_{t+1})$$

- 1-step prediction of the current value:
- 1-step prediction error (called temporal difference (TD) error $\delta_t = r_t + \gamma V(s_{t+1}) V(s_t)$

Update current value with the temporal difference error $V_{\text{new}}(s_t) = V(s_t) + \alpha \delta_t = (1 - \alpha)V(s_t) + \alpha(r_t + \gamma V(s_{t+1}))$



The **TD error**

$$\delta_t = r_t + \gamma V(s_{t+1}) - V(s_t)$$

compares the one-time step lookahead prediction

$$\hat{V}(s_t) = r_t + \gamma V(s_{t+1})$$

with the current estimate of the value function

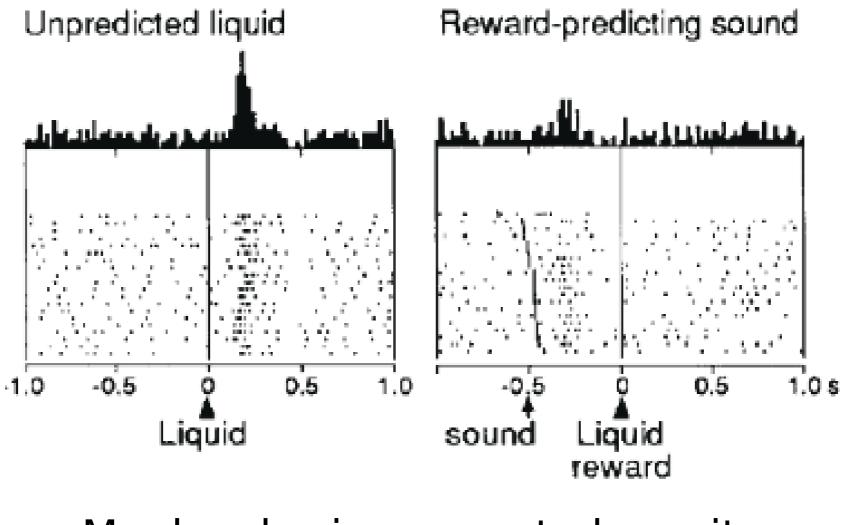
$$\hat{V}(s_t) > V(s_t)$$
 than $V(s_t)$ increased

$$\dot{V}(s_t)$$
 than (s_t) is decreased

Dopamine as TD-error?



Temporal difference error signals can be measured in the brain of monkeys



Monkey brains seem to have it...



Algorithmic Description of TD Learning

 $\mathbf{Init:} V_0^*(s) \leftarrow 0$

Repeat t = t + 1

Observe transition (s_t, a_t, r_t, s_{t+1})

Compute TD error $\delta_t = r_t + \gamma V_t(s_{t+1}) - V_t(s_t)$

Update V-Function $V_{t+1}(s_t) = V_t(s_t) + \alpha \delta_t$

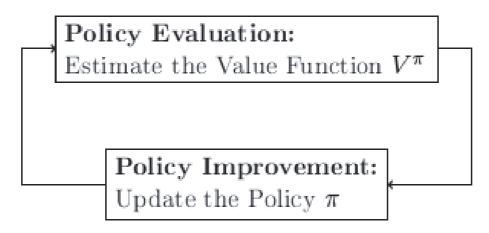
until convergence of V

Used to compute Value function of behavior policySample-based version of policy evaluation



Temporal difference learning for control

So far: Policy evaluation with TD methods



Can we also do the policy improvement step with samples?

Yes, but we need to enforce exploration!

Epsilon-Greedy Policy $(a|s) = \begin{cases} 1 - \epsilon + \epsilon/|\mathcal{A}|, \text{ if } a = \operatorname{argmax}_{a'}Q^{\pi}(s, a') \\ \epsilon/|\mathcal{A}, \text{ otherwise} \end{cases}$

Soft-Max Policy:
$$\pi(\boldsymbol{a}|\boldsymbol{s}) = \frac{\exp(\beta Q(\boldsymbol{s}, \boldsymbol{a}))}{\sum_{\boldsymbol{a}'} \exp(\beta Q(\boldsymbol{s}, \boldsymbol{a}'))}$$

o not always take greedy action

Temporal difference learning for control



Update equations for learning the Q-function

 $Q_{t+1}(s_t, a_t) = Q_t(s_t, a_t) + \alpha \delta_t, \quad \delta_t = r_t + \gamma Q_t(s_{t+1}, a_?) - Q_t(s_t, a_t)$

Two different methods to estimate

Q-learning: $a_? = \operatorname{argmax}_a Q_t(s_{t+1}, a)$

Estimates Q-function of optimal policy

Off-policy samples: $a_? \neq a_{t+1}$

SARSA: $a_? = a_t$, where $a_{t+1} \sim \pi(a|s_{t+1})$

Estimates Q-function of exploration policy

On-policy samples

Note: The policy for generating the actions depends on the Q-function on stationary policy



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In the continuous case, we need to approximate the V-function (except for LQR)

Lets keep it simple, we use a **linear model** to represent the V-function

$$V^{\pi}(\boldsymbol{s}) \approx V_{\boldsymbol{\omega}}(\boldsymbol{s}) = \boldsymbol{\phi}^{T}(\boldsymbol{s})\boldsymbol{\omega}$$

How can we find the parameters ? Again with Temporal Difference Learning



TD-learning with Function Approximation

Derivation:

Use the **recursive definition of V-function**:

$$\begin{aligned} \text{MSE}(\boldsymbol{\omega}) &\approx \text{MSE}_{\text{BS}}(\boldsymbol{\omega}) = 1/N \sum_{i=1}^{N} \left(\hat{V}^{\pi}(\boldsymbol{s}_{i}) - V_{\boldsymbol{\omega}}(\boldsymbol{s}_{i}) \right)^{2} \\ \text{with} \quad \hat{V}^{\pi}(\boldsymbol{s}) = \mathbb{E}_{\pi} \left[r(\boldsymbol{s}, \boldsymbol{a}) + \mathbb{E}_{\mathcal{P}} \left[V_{\boldsymbol{\omega}_{\text{old}}}(\boldsymbol{s}') | \boldsymbol{s}, \boldsymbol{a} \right] \right] \end{aligned}$$

Sootstrapping (BS): Use the old approximation to get the target values for a new approximation

How can we **minimize** this function ?

Lets use **stochastic gradient descent**



Consider an expected error function,

$$E_{\boldsymbol{\omega}} = \mathbb{E}_p[e_{\boldsymbol{\omega}}(x)] \approx 1/N \sum_{i=1}^N e_{\boldsymbol{\omega}}(x_i), \quad x_i \sim p(x)$$

We can find a local minimum of E by Gradient descent:

$$\boldsymbol{\omega}_{k+1} = \boldsymbol{\omega}_k - \alpha_k \frac{dE_{\boldsymbol{\omega}}}{d\boldsymbol{\omega}} = \boldsymbol{\omega}_k - \alpha_k \sum_{i=1}^N \frac{de_{\boldsymbol{\omega}}(x_i)}{d\boldsymbol{\omega}}$$

Stochastic Gradient Descent does the gradient update already after a **single sample**

$$\boldsymbol{\omega}_{k+1} = \boldsymbol{\omega}_k - \alpha_k \frac{de_{\boldsymbol{\omega}}(x_k)}{d\boldsymbol{\omega}}$$

Converges under the stochastic approximation conditions

$$\sum_{k=1}^{\infty} \alpha_k = \infty, \quad \sum_{k=1}^{\infty} \alpha_k^2 < \infty$$



Stochastic gradient descent on our error function *MSE*_{BS}

$$\begin{split} \text{MSE}_{\text{BS,t}}(\boldsymbol{\omega}) &= 1/N \sum_{i=1}^{N} \left(\hat{V}(\boldsymbol{s}_{t}) - V_{\boldsymbol{\omega}}(\boldsymbol{s}_{i}) \right)^{2} \\ &= 1/N \sum_{i=1}^{N} \left(r_{i} + \gamma V_{\boldsymbol{\omega}_{t}}(\boldsymbol{s}_{i}') - V_{\boldsymbol{\omega}}(\boldsymbol{s}_{i}) \right)^{2} \\ \textbf{Update rule (for current time step t, } \boldsymbol{\omega}_{t+1} &= \boldsymbol{\omega}_{t} + \alpha_{t} \left. \frac{d\text{MSE}_{\text{BS}}}{d\boldsymbol{\omega}} \right|_{\boldsymbol{\omega}=\boldsymbol{\omega}_{t}} \\ \boldsymbol{\omega}_{t+1} &= \boldsymbol{\omega}_{t} + \alpha_{t} \left. \frac{d\text{MSE}_{\text{BS}}}{d\boldsymbol{\omega}} \right|_{\boldsymbol{\omega}=\boldsymbol{\omega}_{t}} \\ \boldsymbol{\omega}_{t+1} &= \boldsymbol{\omega}_{t} + \alpha \left(r(\boldsymbol{s}_{t}, \boldsymbol{a}_{t}) + \gamma V_{\boldsymbol{\omega}_{t}}(\boldsymbol{s}_{t+1}) - V_{\boldsymbol{\omega}_{t}}(\boldsymbol{s}_{t}) \right) \boldsymbol{\phi}^{T}(\boldsymbol{s}_{t}) \\ &= \boldsymbol{\omega}_{t} + \alpha \delta_{t} \boldsymbol{\phi}^{T}(\boldsymbol{s}_{t}) \\ \end{split}$$
with
$$\begin{split} \delta_{t} &= r(\boldsymbol{s}_{t}, \boldsymbol{a}_{t}) + \gamma V_{\boldsymbol{\omega}_{t}}(\boldsymbol{s}_{t+1}) - V_{\boldsymbol{\omega}_{t}}(\boldsymbol{s}_{t}) \end{split}$$

with



TD with function approximation

$$\boldsymbol{\omega}_t = \boldsymbol{\omega}_t + \alpha \delta_t \boldsymbol{\phi}^T(\boldsymbol{s}_t)$$

Difference to discrete algorithm:

- TD-error is correlated with the feature vector
- Fixed Equivalent if tabular feature coding is used, $\phi(\mathbf{s}_{i}) = e_{i}$

Similar update rules can be obtained for SARSA and Qlearning $\omega_{t+1} = \omega_t + \alpha \Big(r(s_t, a_t) + \gamma Q_{\omega_t}(s_{t+1}, a_?) - Q_{\omega_t}(s_t, a_t) \Big) \phi^T(s_t, a_t)$

$$Q_{\boldsymbol{\omega}}(\boldsymbol{s}, \boldsymbol{a}) \approx \boldsymbol{\phi}^T(\boldsymbol{s}, \boldsymbol{a}) \boldsymbol{\omega}$$

where

2



Some remarks on temporal difference learning:

- Its not a proper stochastic gradient descent!!
- ➡ Why? Target value $\hat{k}^{\pi}(s)$ change after each parameter update! $\hat{V}^{\pi}(s) = \hat{V}^{\pi}(s)$

We ignore the fact that also depends on

- Side note: This "ignorance" actually introduces a bias in our optimization, such that we are optimizing a different objective than the MSE
- In certain cases, we also get divergence (e.g. off-policy samples)
- TD-learning is very fast in terms of computation time
- 2 O(#features), anbuilteners: data setting leaving with Jeroposal Differences: just sus and comparison, JMLR, in press

Sucessful examples

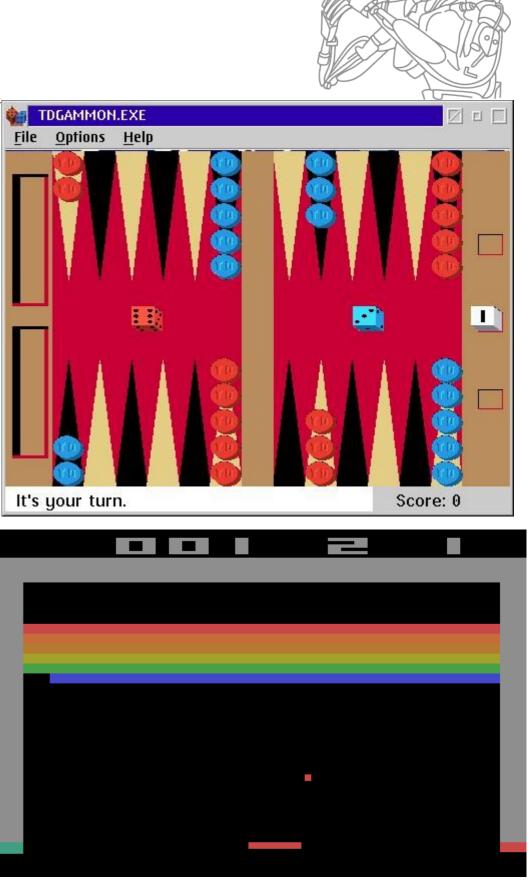
Linear function approximation

Tetris, Go

Non-linear function approximation

TD Gammon (Worldchampion level)

Atari Games (learning from raw pixel input)





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Online methods are typically **data-inefficient** as they use each data point only once

$$D = \left\{ \boldsymbol{s}_i, \boldsymbol{a}_i, r_i, \boldsymbol{s}'_i \right\}_{i=1...N}$$

Can we re-use the whole "batch" of data to increase dataefficiency?

- Least-Squares Temporal Difference (LSTD) Learning
 - **Fitted Q-Iteration**

Computationally much more expensive then TD-learning!



Lets minimize the bootstrapped MSE objective (MSE_{BS})

$$MSE_{BS} = 1/N \sum_{i=1}^{N} \left(r(\boldsymbol{s}_i, \boldsymbol{a}_i) + \gamma V_{\boldsymbol{\omega}_{old}}(\boldsymbol{s}'_i) - V_{\boldsymbol{\omega}}(\boldsymbol{s}_i) \right)^2$$
$$= 1/N \sum_{i=1}^{N} \left(r(\boldsymbol{s}_i, \boldsymbol{a}_i) + \gamma \boldsymbol{\phi}^T(\boldsymbol{s}'_i) \boldsymbol{\omega}_{old} - \boldsymbol{\phi}^T(\boldsymbol{s}_i) \boldsymbol{\omega} \right)^2$$

Least-Squares Solution:

$$\boldsymbol{\omega} = (\boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1} \boldsymbol{\Phi}^T (\boldsymbol{R} + \gamma \boldsymbol{\Phi}' \boldsymbol{\omega}_{\text{old}})$$

with $\mathbf{\Phi} = \left[\boldsymbol{\phi}(\boldsymbol{s}_1), \boldsymbol{\phi}(\boldsymbol{s}_2), \dots, \boldsymbol{\phi}(\boldsymbol{s}_N) \right]^T$ $\mathbf{\Phi}' = \left[\boldsymbol{\phi}(\boldsymbol{s}_1'), \boldsymbol{\phi}(\boldsymbol{s}_2'), \dots, \boldsymbol{\phi}(\boldsymbol{s}_N') \right]^T$

2



Least-Squares Solution:

 $\boldsymbol{\omega} = (\boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1} \boldsymbol{\Phi}^T (\boldsymbol{R} + \gamma \boldsymbol{\Phi}' \boldsymbol{\omega}_{\text{old}})$

Fixed Point: In case of convergence, we want $\omega_{
m old} = \omega$

$$\boldsymbol{\omega} = (\boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1} \boldsymbol{\Phi}^T (\boldsymbol{R} + \gamma \boldsymbol{\Phi}' \boldsymbol{\omega})$$
$$\left(\boldsymbol{I} - \gamma (\boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1} \boldsymbol{\Phi}^T \boldsymbol{\Phi}' \right) \boldsymbol{\omega} = (\boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1} \boldsymbol{\Phi}^T \boldsymbol{R}$$
$$\left(\boldsymbol{\Phi}^T \boldsymbol{\Phi}\right)^{-1} \boldsymbol{\Phi}^T \left(\boldsymbol{\Phi} - \gamma \boldsymbol{\Phi}' \right) \boldsymbol{\omega} = (\boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1} \boldsymbol{\Phi}^T \boldsymbol{R}$$
$$\boldsymbol{\Phi}^T \left(\boldsymbol{\Phi} - \gamma \boldsymbol{\Phi}' \right) \boldsymbol{\omega} = \boldsymbol{\Phi}^T \boldsymbol{R}$$

$$\boldsymbol{\omega} = \left(\boldsymbol{\Phi}^T (\boldsymbol{\Phi} - \gamma \boldsymbol{\Phi}')\right)^{-1} \boldsymbol{\Phi}^T \boldsymbol{R}$$

2



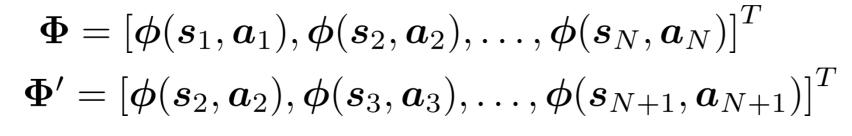
LSTD solution:

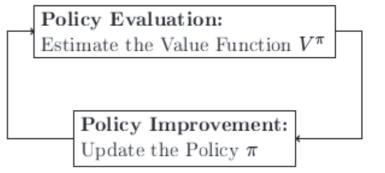
$$\boldsymbol{\omega} = \left(\boldsymbol{\Phi}^T (\boldsymbol{\Phi} - \gamma \boldsymbol{\Phi}') \right)^{-1} \boldsymbol{\Phi}^T \boldsymbol{R}$$

Same solution as convergence point of TD-learning

One shot! No iterations necessary for policy evaluation

LSQ: Adaptation for learning the Q-function





Sed for Least-Squares Policy Iteration (LSPI)

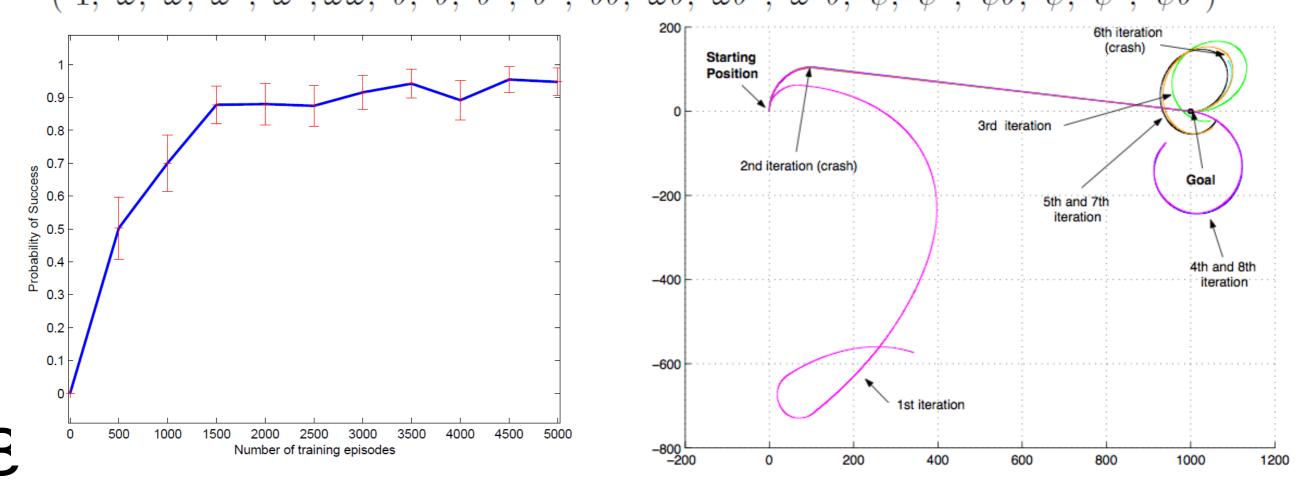
Lagoudakis and Parr, Least-Squares Policy Iteration, JMLR



State space:= $[\theta, \dot{\theta}, \omega, \dot{\omega}, \ddot{\omega}, \psi]$

 θ of handlebar, ω vertical angle of bike, angle to goal **Action space:** 5 discrete actions (torque applied to handle, displacement of rider)

Feature space: 20 basis functions... (1, ω , $\dot{\omega}$, ω^2 , $\dot{\omega}^2$, $\omega\dot{\omega}$, θ , $\dot{\theta}$, θ^2 , $\dot{\theta}^2$, $\theta\dot{\theta}$, $\omega\theta$, $\omega\theta^2$, $\omega^2\theta$, ψ , ψ^2 , $\psi\theta$, $\bar{\psi}$, $\bar{\psi}^2$, $\bar{\psi}\theta$)^T





In Batch-Mode RL it is also much easier to use **non-linear function approximators**

- Many of them only exists in the batch setup, e.g. regression trees
- No catastrophic forgetting, e.g., for neural networks.
- Strong divergence problems, fixed for Neural Networks by ensuring that there is a goal state where the Q-Function value is always zero (see Lange et al. below).

Fitted Q-iteration uses non-linear function approximators for **approximate value iteration.**

Ernst, Geurts and Wehenkel, *Tree-Based Batch Mode Reinforcement Learning, JMLR 2005* ange, Gabel and Riedmiller. *Batch Reinforcement Learning, Reinforcement Learning: State of the Art*



Fitted Q-iteration

Given: Dataset $D = \left\{ m{s}_i, m{a}_i, r_i, m{s}'_i
ight\}_{i=1...N}$

Algorithm:

Initialize $Q^{[0]}(s, a)$ put data:

for k = 1 to L

Generate target values:

Learn new Q-function:

$$\tilde{q}_i^{[k]} = r_i + \gamma \max_{\boldsymbol{a}'} Q^{[k-1]}(\boldsymbol{s}_i', \boldsymbol{a}')$$
$$Q^{[k]}(\boldsymbol{s}, \boldsymbol{a}) \leftarrow \operatorname{Regress}(\boldsymbol{X}, \tilde{\boldsymbol{q}}^{[k]})$$

end

Like Value-Iteration, but we use supervised learning methods to approximate the Q-function at each iteration k

Fitted Q-iteration



Some Remarks:

Regression does the expectation for us

 $Q^{[k]}(\boldsymbol{s}, \boldsymbol{a}) \approx \mathbb{E}_{\mathcal{P}}\left[r(\boldsymbol{s}, \boldsymbol{a}) + \gamma \max_{\boldsymbol{a}'} Q^{[k-1]}(\boldsymbol{s}', \boldsymbol{a}')\right]$

The max operator is still hard to solve for continuous action spaces

For continuous actions, see: Neumann and Peters, Fitted Q-iteration by Advantage weighted regression, NIPS, 2008



Case Study I: Learning Defense

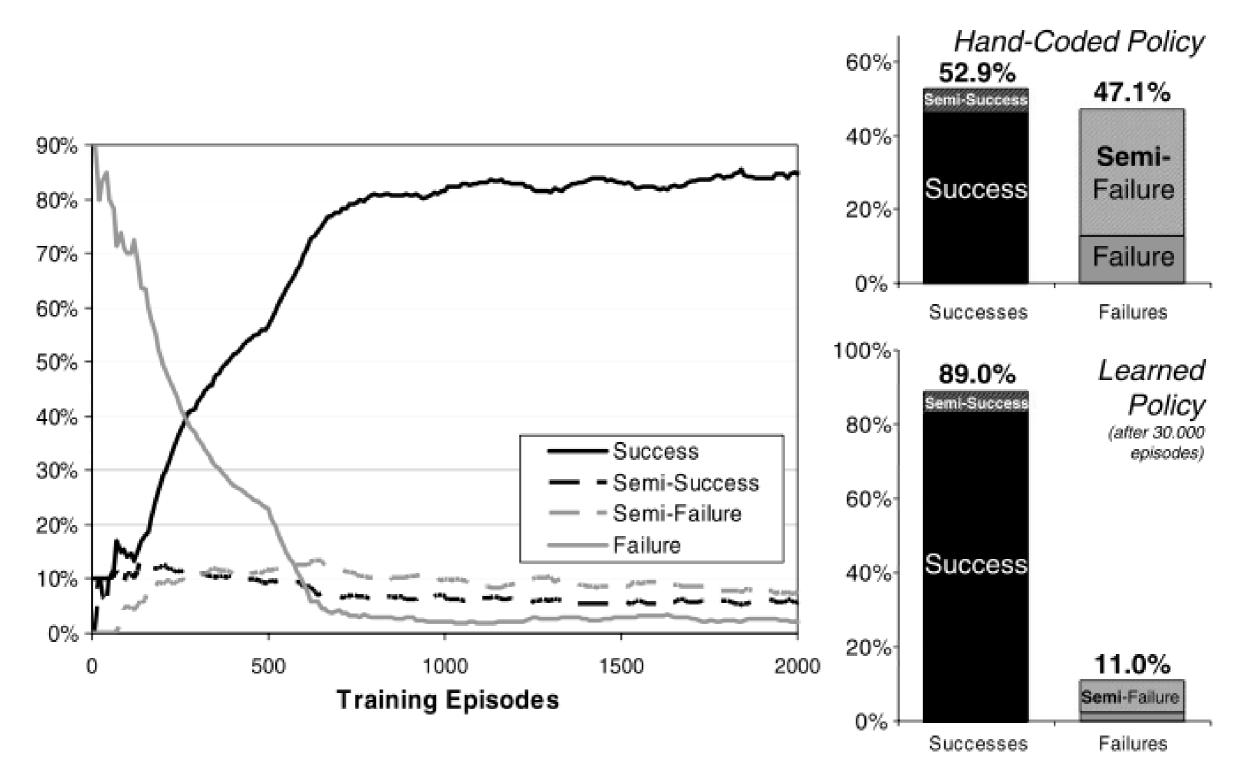


Within the RL framework, we model the ADB learning task as a terminal state problem with both terminal goal S^+ and failure states S^- . Intermediate steps are punished by constant costs of c = 0.05, whereas J(s) = 0.0 for $s \in S^+$ and J(s) = 1.0 for $s \in S^-$ by definition (cf. Eq. 8).

3

Success

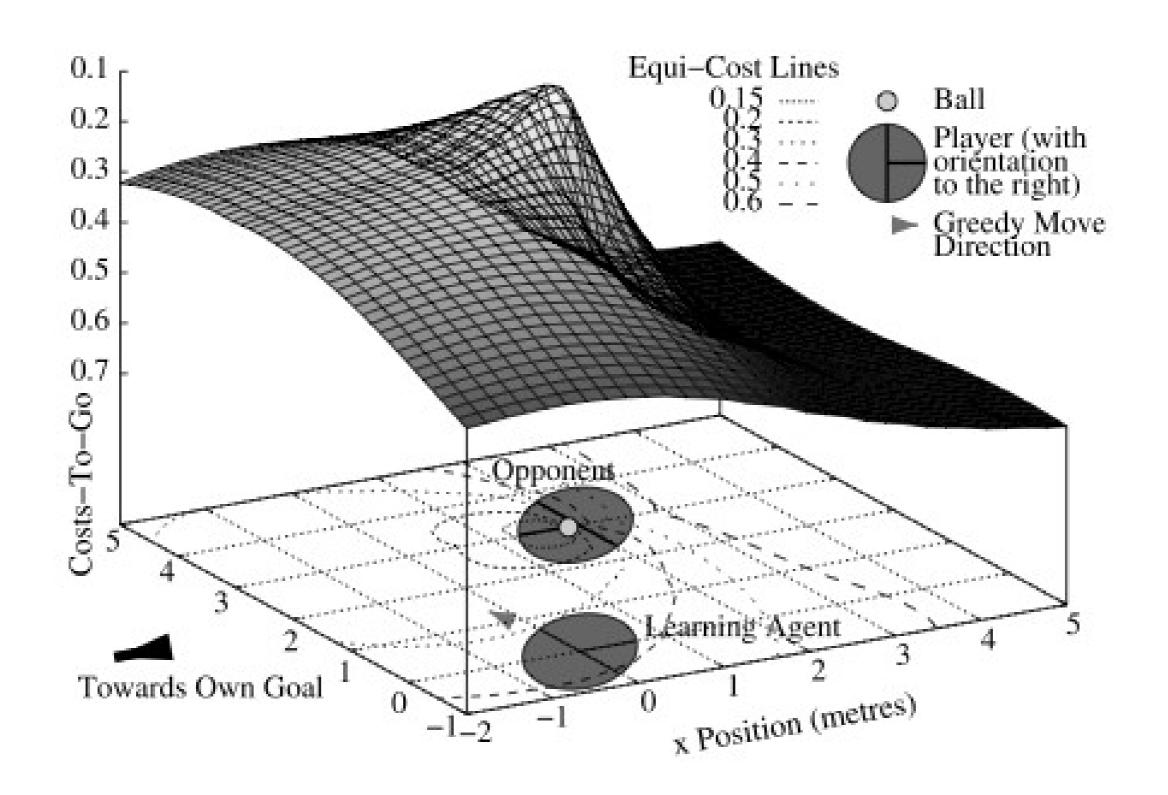




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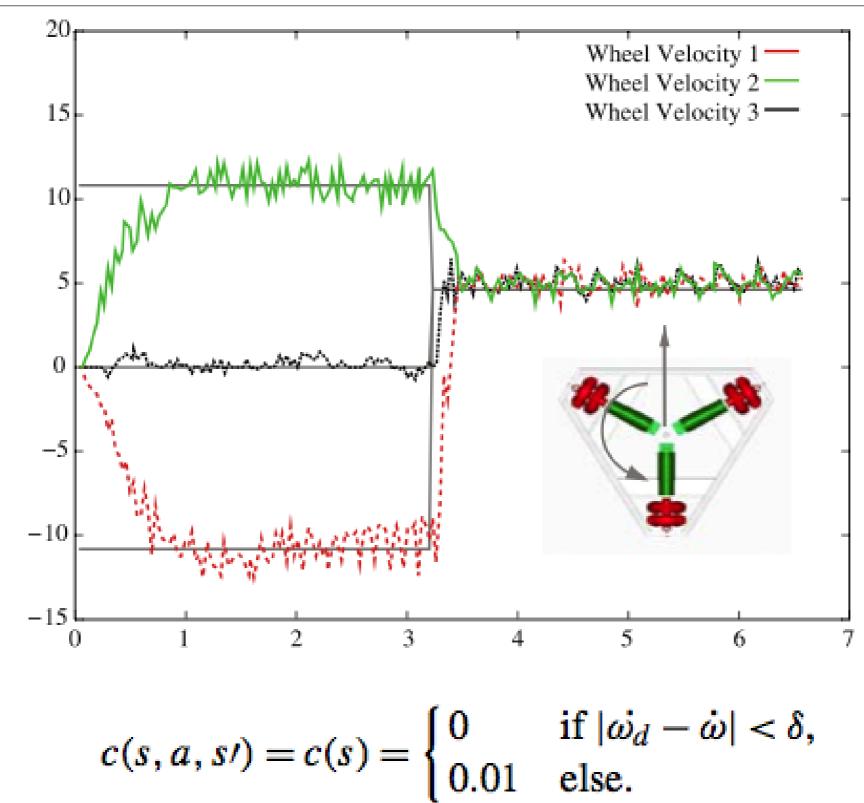
Dueling Behavior





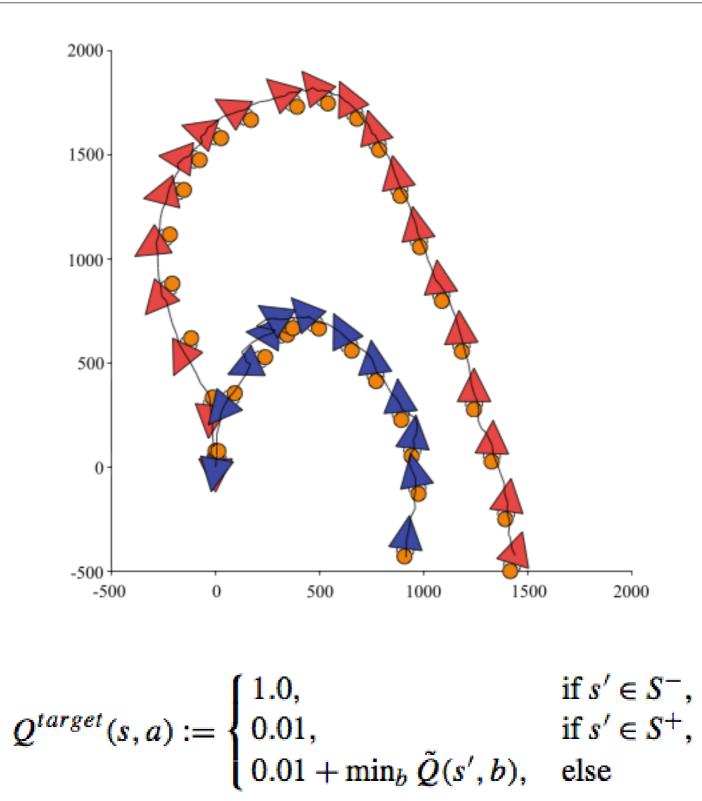


Case Study II: Learning Motor Speeds



3







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- You can do loads of cool things with them: Learn Chess at professional level, learn Backgammon and Checkers at Grandmaster-Level ... and winning the Robot Soccer Cup with a minimum of man power.

So, why are they not always the method of choice?

- You need to fill-up you state-action space up with sufficient samples.
- Another curse of dimensionality with an exponential explosion.
- 3 ➡ Errors in the Value function approximation might have a catastrophic effect on the policy, can be very hard to control