Reinforcement Learning

Value Function Approximation

Stefano V. Albrecht, Michael Herrmann 16 February 2024



- Curse of dimensionality and generalisation
- Value function approximation
- Stochastic gradient descent
- Linear value functions and feature construction
- Semi-gradient TD control

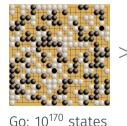
Theory so far has assumed:

- Unlimited space: can store value function as table
- Unlimited data: many (infinite) visits to all state-action pairs

In practice these assumptions are usually violated, because...

Curse of Dimensionality:

- Number of states grows *exponentially* with number of state variables
- If state described by *k* variables with values in {1, ..., *n*}, then *O*(*n^k*) states





Hydrogen atoms: 10⁸⁰

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No data (or not enough data) to estimate return in each state

- Many states may never be visited
- Need to generalise observations to unknown state-action pairs

Generalisation

Blue circle must move to red goal

• Agent uses optimal policy (shortest path)

Suppose we have return estimates (steps to go) for locations S1–S6

• e.g.
$$v(S5) = -3$$
, $v(S4) = -6$, $v(S2) = -31$

We have no data for locations S7 and S8 (not visited yet)

• Can we estimate v(S7) and v(S8) based on other return estimates?

G			S5			S6	
		<mark>S8</mark> ?			S4		
	S1			S2			
	01					<u></u>	
				<mark>S7</mark> ?		S3	

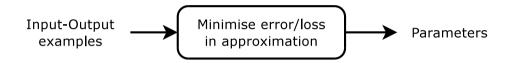
Replace tabular value function with parameterised function:

 $\hat{v}(s, \mathbf{w}) \approx v_{\pi}(s)$ $\hat{q}(s, a, \mathbf{w}) \approx q_{\pi}(s, a)$

 $\mathbf{w} \in \mathbb{R}^d$ is parameter ("weight") vector e.g. linear function, neural network, regression tree, ...

- Compact: number of parameters d much smaller than |S|
- **Generalises:** changing one parameter value may change value estimate of many states/actions

Learning a value function is a form of supervised learning:



Examples are pairs of states and return estimates, (S_t, U_t) , e.g.

- MC: $U_t = G_t$
- TD(0): $U_t = R_{t+1} + \gamma \hat{v}(S_{t+1}, \mathbf{w}_t)$
- n-step TD: $U_t = R_{t+1} + \cdots + \gamma^{n-1}R_{t+n} + \gamma^n \hat{v}(S_{t+n}, \mathbf{w}_{t+n-1})$

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• Incremental updates

update w using only partial data, e.g. most recent (S_t, U_t) or subset

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• Ability to handle noisy targets e.g. different MC updates *G_t* for same state *S_t* Desired properties in supervised learning method:

• Incremental updates

update w using only partial data, e.g. most recent (S_t, U_t) or subset

- Ability to handle noisy targets e.g. different MC updates *G_t* for same state *S_t*
- Ability to handle non-stationary targets e.g. changing target policy, bootstrapping

 \Rightarrow If \hat{v}/\hat{q} differentiable, stochastic gradient descent is suitable method

Gradient Descent

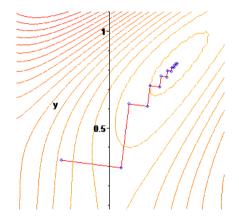
- Let J(w) be differentiable function of w
- Gradient of J(w) is

$$abla J(\mathbf{w}) = \left(\frac{\partial J(\mathbf{w})}{\partial w_1}, \cdots, \frac{\partial J(\mathbf{w})}{\partial w_d}\right)^{\top}$$

• To find local minimum of J(w), adjust w in negative direction of gradient

$$\mathbf{w}_{t+1} = \mathbf{w}_t - \frac{1}{2} \alpha \, \nabla J(\mathbf{w}_t)$$

 α is step-size parameter convergence requires standard α-reduction



Objective: find parameter vector **w** by minimising *mean-squared error* between approximate value $\hat{v}(s, \mathbf{w})$ and true value $v_{\pi}(s)$

$$J(\mathbf{w}) = \mathbb{E}_{\pi} \big[(v_{\pi}(s) - \hat{v}(s, \mathbf{w}))^2 \big]$$

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• Gradient descent finds local minimum:

$$\mathbf{w}_{t+1} = \mathbf{w}_t - \frac{1}{2} \alpha \nabla J(\mathbf{w}_t)$$

= $\mathbf{w}_t + \alpha \mathbb{E}_{\pi}[(\mathbf{v}_{\pi}(s) - \hat{\mathbf{v}}(s, \mathbf{w}_t)) \nabla \hat{\mathbf{v}}(s, \mathbf{w}_t)]$

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• Stochastic gradient descent samples the gradient:

$$\mathbf{w}_{t+1} = \mathbf{w}_t + \alpha \left[\mathbf{U}_t - \hat{\mathbf{v}}(\mathbf{S}_t, \mathbf{w}_t) \right] \nabla \hat{\mathbf{v}}(\mathbf{S}_t, \mathbf{w}_t)$$

Stochastic Gradient Descent – Convergence

Stochastic gradient descent *samples* the gradient:

$$\mathbf{N}_{t+1} = \mathbf{W}_t + \alpha \left[U_t - \hat{\mathbf{v}}(\mathbf{S}_t, \mathbf{W}_t) \right] \nabla \hat{\mathbf{v}}(\mathbf{S}_t, \mathbf{W}_t) \tag{1}$$

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• \mathbf{w}_t will converge to local optimum under standard α -reduction and if U_t is unbiased estimate $\mathbb{E}_{\pi}[U_t|S_t] = v_{\pi}(S_t)$

 \Rightarrow MC update is unbiased but TD update is biased (why?)

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• Note: (1) is not a true TD gradient because U_t also depends on **w**

$$U_t = R_{t+1} + \gamma \hat{v}(S_{t+1}, \mathbf{w})$$

Hence, we call it semi-gradient TD

Input: the policy π to be evaluated Input: a differentiable function $\hat{v}: \mathbb{S}^+ \times \mathbb{R}^d \to \mathbb{R}$ such that $\hat{v}(\text{terminal}, \cdot) = 0$ Algorithm parameter: step size $\alpha > 0$ Initialize value-function weights $\mathbf{w} \in \mathbb{R}^d$ arbitrarily (e.g., $\mathbf{w} = \mathbf{0}$)

Loop for each episode:

```
Initialize S

Loop for each step of episode:

Choose A \sim \pi(\cdot|S)

Take action A, observe R, S'

\mathbf{w} \leftarrow \mathbf{w} + \alpha [R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})] \nabla \hat{v}(S, \mathbf{w})

S \leftarrow S'

until S is terminal
```

Linear Value Function Approximation

Linear value function approximation:

$$\hat{v}(s, \mathbf{w}) \doteq \mathbf{w}^{\top} \mathbf{x}(s) = \sum_{i=1}^{d} w_i x_i(s)$$

- $\mathbf{x}(s) = (x_1(s), ..., x_d(s))^\top$ is feature vector of state s
- Simple gradient: $\nabla \hat{v}(s, \mathbf{w}) = \left(\frac{\partial \mathbf{w}^{\top} \mathbf{x}}{\partial w_1}, \cdots, \frac{\partial \mathbf{w}^{\top} \mathbf{x}}{\partial w_d}\right)^{\top} = \mathbf{x}(s)$
- Gradient update: $\mathbf{w}_{t+1} = \mathbf{w}_t + \alpha \left[U_t \hat{v}(S_t, \mathbf{w}_t) \right] \mathbf{x}(S_t)$

See Tutorial 5

Linear Value Function Approximation

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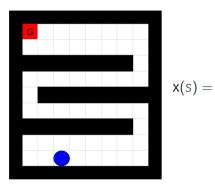
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In linear case, there is only one optimum!

- \Rightarrow MC gradient updates converge to global optimum
- \Rightarrow TD gradient updates converge *near* global optimum (TD fixed point)

See Tutorial 5

Feature Vectors



Remember: State must be Markov

$$\begin{pmatrix} x-\text{pos}(s) \\ y-\text{pos}(s) \end{pmatrix}$$
$$\mathbf{x}(s) = \begin{pmatrix} \theta(s) \\ \theta-\text{vel}(s) \\ x-\text{pos}(s) \\ \vdots \end{pmatrix} \xrightarrow{\mathbf{y}}_{M} \underbrace{\vec{F}}_{M}$$

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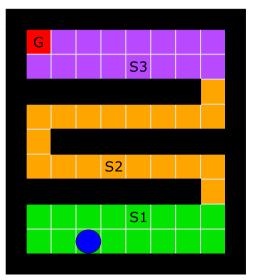
Exact representation:

$$\mathbf{x}(s) = \left(\begin{array}{c} x \text{-} pos(s) \\ y \text{-} pos(s) \end{array}\right)$$

Generalise with state aggregation:

• Partition states into disjoint sets $S_1, S_2, ...$ with indicator functions $\mathbf{x}_k(s) = [s \in S_k]_1$

$$\mathbf{x}(s) = \left(\begin{array}{c} \text{in-S1}(s)\\ \text{in-S2}(s)\\ \text{in-S3}(s) \end{array}\right) = \left(\begin{array}{c} 1\\ 0\\ 0 \end{array}\right)$$



Exact representation:

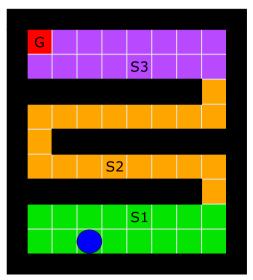
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Generalise with state aggregation:

• Partition states into disjoint sets $S_1, S_2, ...$ with indicator functions $\mathbf{x}_k(s) = [s \in S_k]_1$

Special case: every state s has its own set $\mathcal{S}_s = \{s\}$

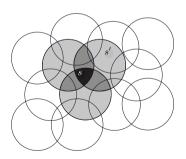
 \Rightarrow Same as tabular representation!

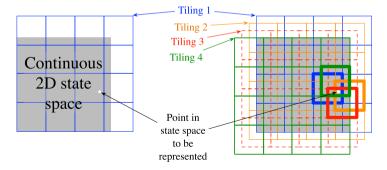


Coarse/Tile Coding

State aggregation generalises only within sets S_1 , S_2 , ...

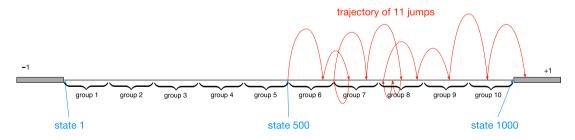
- Allow generalisation *across* sets by allowing S_k to overlap
- e.g. coarse coding and tile coding



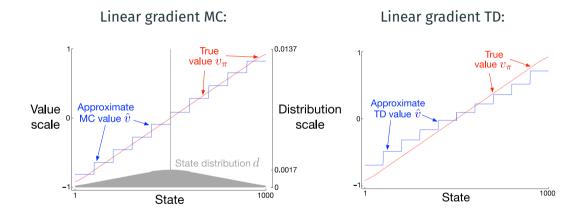


Example: Random Walk

- States: numbered 1 to 1000, start at state 500
- Policy: randomly jump to one of 100 states to left, or one of 100 states to right
- If jump goes beyond 1/1000, terminates with reward -1/+1
- State aggregation: 10 groups of 100 states each



Random Walk: MC and TD Prediction



After 100,000 episodes with $\alpha = 2 \times 10^{-5}$

Approximate Control in Episodic Tasks

- Estimate state-action values: $\hat{q}(s, a, \mathbf{w}) \approx q_{\pi}(s, a)$
- For linear approx., features defined over states and action:

$$\hat{q}(s, a, \mathbf{w}) \doteq \sum_{i=1}^{d} w_i x_i(s, a)$$

• Stochastic gradient descent:

$$\mathbf{w}_{t+1} = \mathbf{w}_t + \alpha \left[U_t - \hat{q}(S_t, A_t, \mathbf{w}_t) \right] \nabla \hat{q}(S_t, A_t, \mathbf{w}_t)$$

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e.g. Sarsa:
$$U_t = R_{t+1} + \gamma \,\hat{q}(S_{t+1}, A_{t+1}, \mathbf{w}_t)$$

Q-learning: $U_t = R_{t+1} + \gamma \max_a \hat{q}(S_{t+1}, a, \mathbf{w}_t)$

Expected Sarsa: $U_t = R_{t+1} + \gamma \sum_a \pi(a|S_{t+1}) \hat{q}(S_{t+1}, a, \mathbf{w}_t)$

Episodic Semi-gradient Sarsa

Input: a differentiable action-value function parameterization $\hat{q} : S \times \mathcal{A} \times \mathbb{R}^d \to \mathbb{R}$ Algorithm parameters: step size $\alpha > 0$, small $\varepsilon > 0$ Initialize value-function weights $\mathbf{w} \in \mathbb{R}^d$ arbitrarily (e.g., $\mathbf{w} = \mathbf{0}$)

Loop for each episode:

 $S, A \leftarrow \text{initial state and action of episode (e.g., <math>\varepsilon$ -greedy)

Loop for each step of episode:

Take action A, observe R, S'

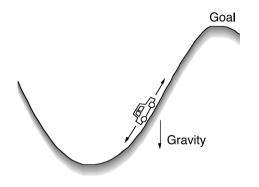
If S' is terminal:

 $\mathbf{w} \leftarrow \mathbf{w} + \alpha \big[R - \hat{q}(S, A, \mathbf{w}) \big] \nabla \hat{q}(S, A, \mathbf{w})$

Go to next episode

Choose A' as a function of $\hat{q}(S', \cdot, \mathbf{w})$ (e.g., ε -greedy) $\mathbf{w} \leftarrow \mathbf{w} + \alpha [R + \gamma \hat{q}(S', A', \mathbf{w}) - \hat{q}(S, A, \mathbf{w})] \nabla \hat{q}(S, A, \mathbf{w})$ $S \leftarrow S'$ $A \leftarrow A'$

Example: Mountain Car with Linear Semi-Gradient Sarsa



STATES: car's position and velocity

<u>ACTIONS</u>: three thrusts: forward, reverse, none

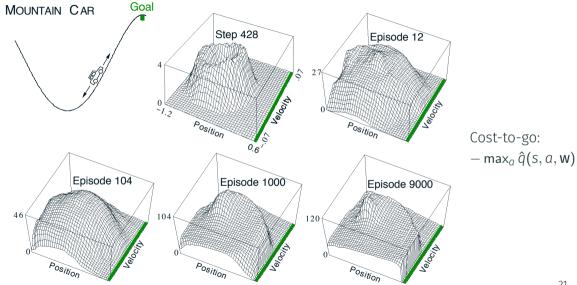
REWARDS:

always -1 until car reaches the goal

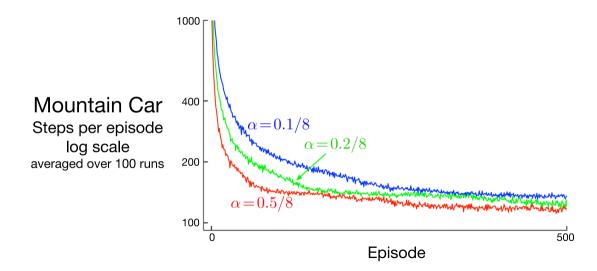
Episodic, No Discounting, $\gamma=1$

Semi-gradient Sarsa with linear approximation over 8 8x8 tilings $\epsilon = 0$ (optimistic initial values $\hat{q}(s, a, \mathbf{w}) = 0$)

Learned Action Values in Mountain Car



Learning Curves in Mountain Car



Algorithm	Tabular	Linear	Non-linear
MC control	yes	chatter*	no
(semi-gradient) n-step Sarsa	yes	chatter*	no
(semi-gradient) n-step Q-learning	yes	no	no

*Chatters near optimal solution because optimal policy may not be representable under value function approximation

Risk of divergence arises when the following three are combined:

- 1. Function approximation
- 2. Bootstrapping
- 3. Off-policy learning

Possible fixes:

- Use importance sampling to warp off-policy distribution into on-policy distribution
- Use gradient TD methods which follow true gradient of projected Bellman error (see book)

Reading

Required (RL book):

• Chapter 9 (9.1–9.5)

(Box "Proof of Convergence of Linear TD(0)" in Sec 9.4 is not examined)

- Chapter 10 (10.1)
- Chapter 11 (11.1)

Optional:

- Remaining sections of chapters
- Tsitsiklis, J. N., Van Roy, B. (1997). An analysis of temporal-difference learning with function approximation. IEEE Transactions on Automatic Control, 42(5):674–690
- Mahadevan, S. (1996). Average reward reinforcement learning: Foundations, algorithms, and empirical results. Machine Learning, 22(1):159–196