

Basics of reinforcement learning

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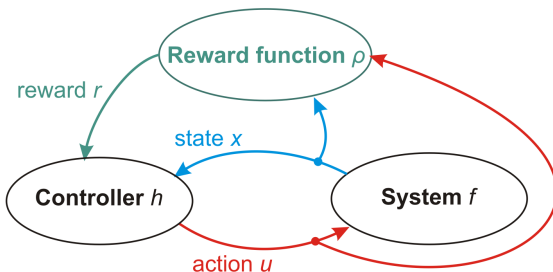
TMLSS, 20 July 2018



Learn a sequential decision policy
to optimize the cumulative performance
of an unknown system

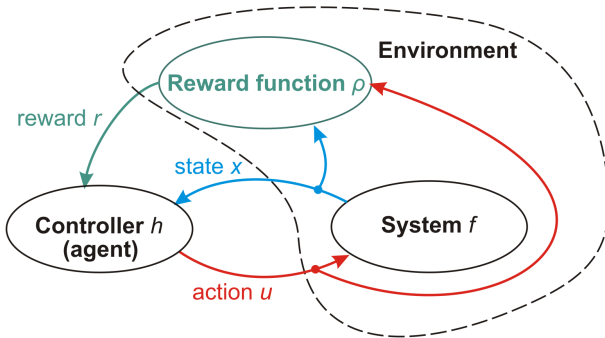


RL principle



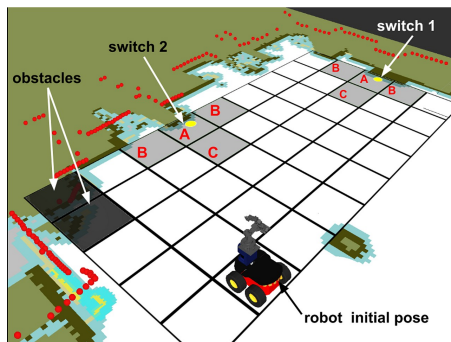
- Interact with system: measure **states**, apply **actions**
- Performance feedback in the form of **rewards**
- Inspired by human and animal learning

RL principle: AI view



- Agent embedded in an environment that feeds back states and rewards

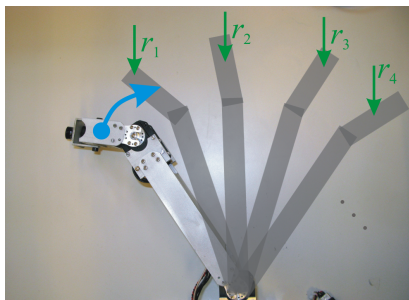
Example: Domestic robot



A domestic robot ensures light switches are off
Abstractization to high-level control (physical actions implemented by low-level controllers)

- **States:** grid coordinates, switch states
- **Actions:** movements NSEW, toggling switch
- **Rewards:** when switches toggled on→off

Example: Robot arm

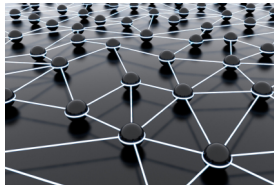
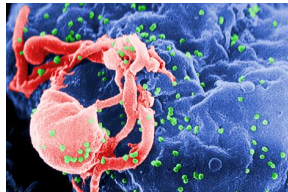
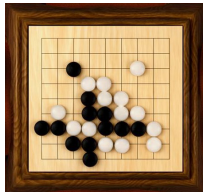


Low-level control

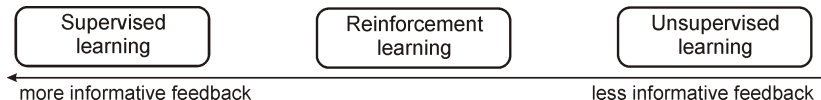
- **States**: link angles and angular velocities
- **Actions**: motor voltages
- **Rewards**: e.g. to reach a desired configuration, give larger rewards as robot gets closer to it

Many other applications

Artificial intelligence, control, medicine, multiagent systems, economics etc.



RL on the machine learning spectrum

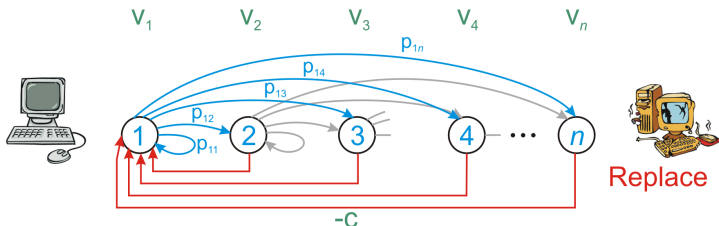


- Supervised: for each training sample, **correct output** known
- Unsupervised: only input samples, **no outputs**; find patterns in the data
- Reinforcement: correct actions not available, **only rewards**

But note: RL finds **dynamical optimal control**!

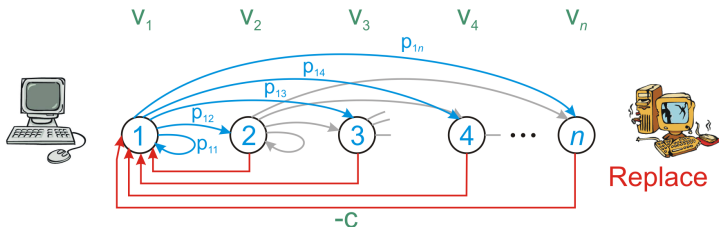
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Example: Machine replacement



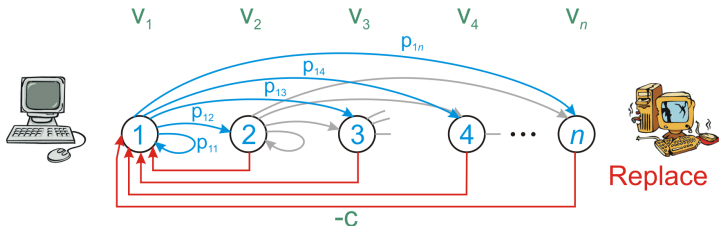
- Consider machine with n different wear levels
1 = perfect working order, n = fully degraded
- Produces revenue v_i when operating at level i
- Stochastic wear: level i increases to $j > i$ with probas p_{ij} ,
stays i with $p_{ii} = 1 - p_{i,i+1} - \dots - p_{i,n}$
- Machine can be replaced at any time (assumed instantaneously), paying cost c

Machine replacement: States and actions



- **State** x = wear level,
state space $X = \{1, 2, \dots, n\}$
- **Action** u = whether to **W**ait or **R**eplace,
action space: $U = \{W, R\}$

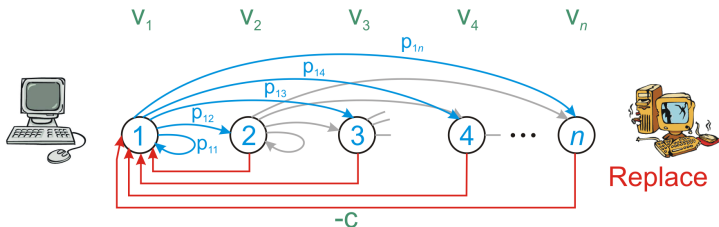
Machine replacement: Transition function



Transition function $f(x, u, x')$ gives the probability of reaching state x' after applying action u in state x :

$$f(x = i, u, x' = j) = \begin{cases} p_{ij} & \text{if } u = W \text{ and } i \leq j \\ 1 & \text{if } u = R \text{ and } j = 1 \\ 0 & \text{otherwise} \end{cases}$$

Machine replacement: Reward function



Reward function $\rho(x, u, x')$ gives reward resulting from transitioning from x to x' after applying u :

$$\rho(x = i, u, x' = j) = \begin{cases} v_j & \text{if } u = W \\ -c + v_1 & \text{if } u = R \end{cases}$$

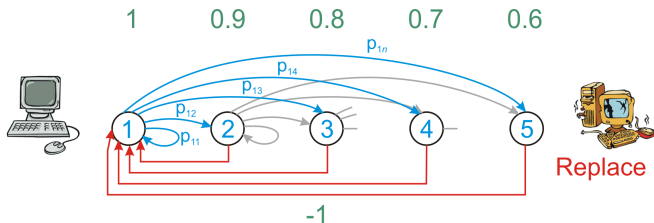
General case: Markov decision process

Markov decision process (MDP)

- 1 State space X
- 2 Action space U
- 3 Transition function $f(x, u, x')$, $f : X \times U \times X \rightarrow [0, 1]$
- 4 Reward function $\rho(x, u, x')$, $\rho : X \times U \times X \rightarrow \mathbb{R}$

Some MDPs have terminal states (e.g. success, failure), that once reached cannot be left and provide no additional reward

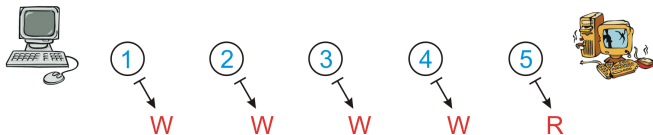
Machine replacement: Specific example



- $n = 5$ wear levels
- Revenue: $v_1 = 1, v_2 = 0.9, \dots, v_5 = 0.6$
- Cost of new machine: $c = 1$
- Wear increase probabilities:

$$[p_{ij}] = \begin{bmatrix} 0.6 & 0.3 & 0.1 & 0 & 0 \\ 0 & 0.6 & 0.3 & 0.1 & 0 \\ 0 & 0 & 0.6 & 0.3 & 0.1 \\ 0 & 0 & 0 & 0.7 & 0.3 \\ 0 & 0 & 0 & 0 & 1.0 \end{bmatrix}$$

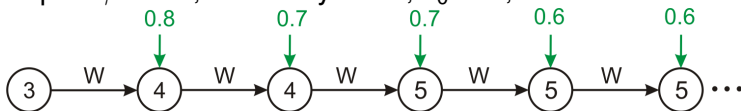
Control policy



- **Control policy** $h : X \rightarrow U$: maps states x to actions u
- Example for machine replacement: $h(1) = \dots = h(4) = W$, $h(5) = R$

Return and objective

Example: $\gamma = 0.9$, $h = \text{always wait}$, $x_0 = 4$, and trial:



$$\gamma^0 r_1 + \gamma^1 r_2 + \gamma^2 r_3 + \gamma^3 r_4 + \dots = 0.9^0 \cdot 0.8 + 0.9^1 \cdot 0.7 + \\ + 0.9^2 \cdot 0.7 + 0.9^3 \cdot 0.6 + 0.9^4 \cdot 0.6 + \dots = 6.3710$$

Objective

Find h that maximizes from any x_0
the **expected return** under the stochastic transitions:

$$R^h(x_0) = E_{x_1, x_2, \dots} \left\{ \sum_{k=0}^{\infty} \gamma^k \rho(x_k, h(x_k), x_{k+1}) \right\}$$

Note: There are also other types of return!



Discount factor

Discount factor $\gamma \in [0, 1)$:

- represents an increasing uncertainty about the future
- bounds the infinite sum (assuming rewards bounded)
- helps the convergence of algorithms

To choose γ , **trade-off** between:

- 1 Long-term quality of the solution (large γ)
- 2 Simplicity of the problem (small γ)

In practice, γ should be sufficiently large so as not to ignore important later rewards



Q-function

Q-function of a policy h is the expected return achieved by executing u_0 in x_0 and then following h

$$Q^h(x_0, u_0) = E_{x_1} \left\{ \rho(x_0, u_0, x_1) + \gamma R^h(x_1) \right\}$$

Q^h measures the quality of state-action pairs



Bellman equation of Q^h

Go one step further in the equation:

$$\begin{aligned} Q^h(x_0, u_0) &= E_{x_1} \left\{ \rho(x_0, u_0, x_1) + \gamma R^h(x_1) \right\} \\ &= E_{x_1} \left\{ \rho(x_0, u_0, x_1) + \gamma E_{x_2} \left\{ \rho(x_1, h(x_1), x_2) + \gamma R^h(x_2) \right\} \right\} \\ &= E_{x_1} \left\{ \rho(x_0, u_0, x_1) + \gamma Q^h(x_1, h(x_1)) \right\} \end{aligned}$$

⇒ **Bellman equation for Q^h**

$$\begin{aligned} Q^h(x, u) &= E_{x'} \left\{ \rho(x, u, x') + \gamma Q^h(x', h(x')) \right\} \\ &= \sum_{x'} f(x, u, x') \left[\rho(x, u, x') + \gamma Q^h(x', h(x')) \right] \end{aligned}$$

Optimal solution and Bellman optimality equation

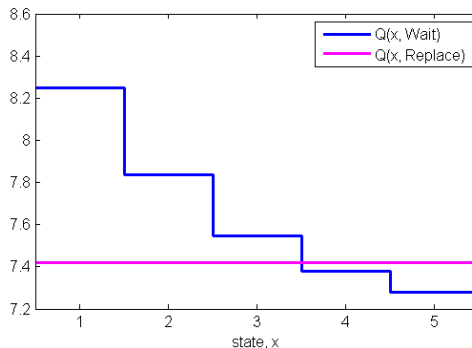
- **Optimal Q-function:** $Q^* = \max_h Q^h$
- ⇒ “Greedy” policy in Q^* : $h^*(x) = \arg \max_u Q^*(x, u)$
 is **optimal**, i.e. achieves maximal returns
 (if multiple actions maximize, break ties arbitrarily)

Bellman optimality equation (for Q^*)

$$\begin{aligned}
 Q^*(x, u) &= E_{x'} \left\{ \rho(x, u, x') + \gamma \max_{u'} Q^*(x', u') \right\} \\
 &= \sum_{x'} f(x, u, x') \left[\rho(x, u, x') + \gamma \max_{u'} Q^*(x', u') \right]
 \end{aligned}$$

Machine replacement: Optimal solution

Discount factor $\gamma = 0.9$



Up next:

Algorithms to find the optimal solution



Algorithm landscape

By model usage:

- **Model-based**: f, ρ known a priori
- **Model-free**: f, ρ unknown
- **Model-learning**: f, ρ learned from data

Model-based usually called dynamic programming (DP);
needed as a stepping stone to RL

By interaction level:

- **Offline**: algorithm runs in advance
- **Online**: algorithm runs with the system

We focus on **exact** case: x, u small number of discrete values,
so we can exactly represent solutions. In practice, function
approximation often needed – covered in Doina's talk



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Q-iteration

Transforms Bellman optimality equation:

$$Q^*(x, u) = \sum_{x'} f(x, u, x') \left[\rho(x, u, x') + \gamma \max_{u'} Q^*(x', u') \right]$$

into an **iterative procedure**:

Q-iteration

repeat at each iteration ℓ

for all x, u **do**

$$Q_{\ell+1}(x, u) \leftarrow \sum_{x'} f(x, u, x') \left[\rho(x, u, x') + \gamma \max_{u'} Q_{\ell}(x', u') \right]$$

end for

until convergence to Q^*

Once Q^* available: $h^*(x) = \arg \max_u Q^*(x, u)$

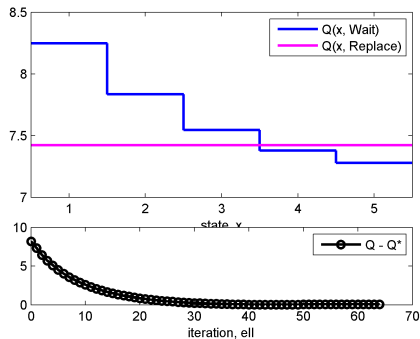
Q-iteration belongs to the class of value iteration algorithms



Machine replacement: Q-iteration demo

Discount factor $\gamma = 0.9$

Q-iteration, $\text{ell}=64$



Machine replacement: Q-iteration demo

$$Q_{\ell+1}(x, u) \leftarrow \sum_{x'} f(x, u, x') [\rho(x, u, x') + \gamma \max_{u'} Q_{\ell}(x', u')]$$

	$x = 1$	$x = 2$	$x = 3$	$x = 4$	$x = 5$
Q_0	0 ; 0	0 ; 0	0 ; 0	0 ; 0	0 ; 0
Q_1	1 ; 0	0.9 ; 0	0.8 ; 0	0.7 ; 0	0.6 ; 0
Q_2	1.86 ; 0.9	1.67 ; 0.9	1.48 ; 0.9	1.3 ; 0.9	1.14 ; 0.9
Q_3	2.58 ; 1.67	2.31 ; 1.67	2.05 ; 1.67	1.83 ; 1.67	1.63 ; 1.67
Q_4	3.2 ; 2.33	2.87 ; 2.33	2.55 ; 2.33	2.3 ; 2.33	2.1 ; 2.33
...
Q_{64}	8.25 ; 7.42	7.84 ; 7.42	7.55 ; 7.42	7.38 ; 7.42	7.28 ; 7.42
Q_{65}	8.25 ; 7.42	7.84 ; 7.42	7.55 ; 7.42	7.38 ; 7.42	7.28 ; 7.42
h^*	W	W	W	R	R

$$h^*(x) = \arg \max_u Q^*(x, u)$$



Policy iteration

Policy iteration

initialize policy h_0

repeat at each iteration ℓ

1: **policy evaluation**: find Q^{h_ℓ}

2: **policy improvement**:

$$h_{\ell+1}(x) \leftarrow \arg \max_u Q^{h_\ell}(x, u)$$

until convergence to h^*

Iterative policy evaluation

Similarly to Q-iteration, transforms Bellman equation for Q^h :

$$Q^h(x, u) = \sum_{x'} f(x, u, x') \left[\rho(x, u, x') + \gamma Q^h(x', h(x')) \right]$$

into an iterative procedure:

Iterative policy evaluation

repeat at each iteration τ

for all x, u **do**

$$Q_{\tau+1}(x, u) \leftarrow \sum_{x'} f(x, u, x') \left[\rho(x, u, x') + \gamma Q_{\tau}(x', h(x')) \right]$$

end for

until convergence to Q^h

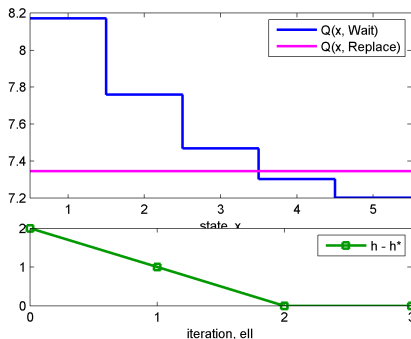
(other options exist, e.g. solving the linear system)



Machine replacement: policy iteration demo

Discount factor $\gamma = 0.9$

Policy iteration, $\text{ell}=3$



Machine replacement: policy iteration

$$Q_{\tau+1}(x, u) \leftarrow \sum_{x'} f(x, u, x') [\rho(x, u, x') + \gamma Q_{\tau}(x', h(x'))]$$

$$h_{\ell+1}(x) \leftarrow \arg \max_u Q^{h_{\ell}}(x, u)$$

	$x = 1$	$x = 2$	$x = 3$	$x = 4$	$x = 5$
h_0	W	W	W	W	W
Q_0	0; 0	0; 0	0; 0	0; 0	0; 0
Q_1	1; 0	0.9; 0	0.8; 0	0.7; 0	0.6; 0
Q_2	1.86; 0.9	1.67; 0.9	1.48; 0.9	1.3; 0.9	1.14; 0.9
Q_3	2.58; 1.67	2.31; 1.67	2.05; 1.67	1.83; 1.67	1.63; 1.67
...
Q_{39}	7.51; 6.75	6.95; 6.75	6.49; 6.75	6.17; 6.75	5.9; 6.75
Q_{40}	7.52; 6.75	6.96; 6.75	6.5; 6.75	6.18; 6.75	5.91; 6.75
h_1	W	W	R	R	R

...algorithm continues...



Machine replacement: policy iteration (cont'd)

	$x = 1$	$x = 2$	$x = 3$	$x = 4$	$x = 5$
h_1	W	W	R	R	R
Q_0	0; 0	0; 0	0; 0	0; 0	0; 0
...
Q_{43}	8.01; 7.2	7.57; 7.2	7.27; 7.2	7.17; 7.2	7.07; 7.2
h_2	W	W	W	R	R
Q_0	0; 0	0; 0	0; 0	0; 0	0; 0
...
Q_{43}	8.17; 7.35	7.76; 7.35	7.47; 7.35	7.3; 7.35	7.2; 7.35
h_3	W	W	W	R	R

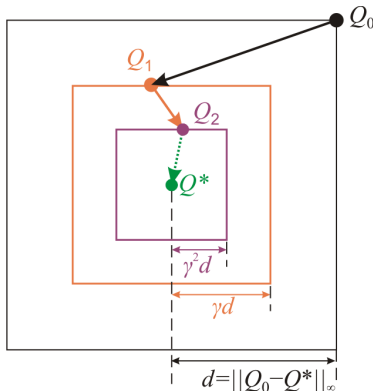


Convergence of Q-iteration

- Each iteration a contraction with factor γ in ∞ -norm:

$$\|Q_{\ell+1} - Q^*\|_{\infty} \leq \gamma \|Q_{\ell} - Q^*\|_{\infty}$$

⇒ Q-iteration **monotonically converges** to Q^* ,
with convergence rate $\gamma \Rightarrow \gamma$ helps convergence



Stopping condition of Q-iteration

- Convergence to Q^* only guaranteed asymptotically, as $\ell \rightarrow \infty$
- In practice, algorithm can be stopped when:

$$\|Q_{\ell+1} - Q_{\ell}\| \leq \varepsilon_{\text{qiter}}$$



Convergence of policy iteration

Policy evaluation component – like Q-iteration:

- Iterative policy evaluation contraction with factor γ
- ⇒ **monotonic convergence** to Q^h , with rate γ

Complete policy iteration algorithm:

- Policy is either improved or already optimal
 - But the maximum number of improvements is finite! ($|U|^{|X|}$)
- ⇒ **convergence** to h^* in a finite number of iterations



Stopping conditions of policy iteration

In practice:

- Policy evaluation can be stopped when:

$$\|Q_{\tau+1} - Q_{\tau}\| \leq \varepsilon_{\text{peval}}$$

- Policy iteration can be stopped when:

$$\|h_{\ell+1} - h_{\ell}\| \leq \varepsilon_{\text{piter}}$$

- Note: $\varepsilon_{\text{piter}}$ can be taken 0!

Q-iteration vs. policy iteration

Number of iterations to convergence

- Q-iteration $>$ policy iteration

Complexity

- one iteration of Q-iteration
 $>$ one iteration of iterative policy evaluation
- complete Q-iteration ??? complete policy iteration



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By interaction level:

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We move to **online RL** for the remainder of the talk



Policy evaluation change

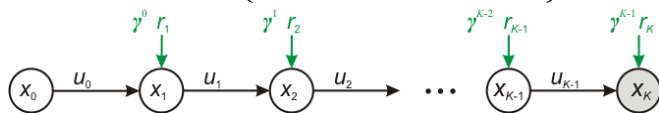
To find Q^h :

- So far: model-based policy evaluation
- Reinforcement learning: model not available!
- Learn Q^h from data obtained by
online interaction with the system



Monte Carlo policy evaluation

Recall: $Q^h(x_0, u_0) = E_{x_1} \{ \rho(x_0, u_0, x_1) + \gamma R^h(x_1) \}$



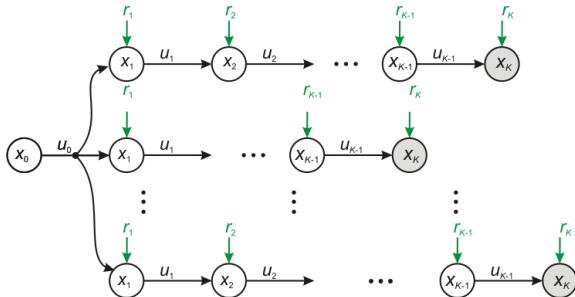
- Trial from (x_0, u_0) to x_K using $u_1 = h(x_1)$, $u_2 = h(x_2)$, etc.
 - x_K must be terminal (assumed further) or K large enough
- ⇒ Return along trial provides a sample of $Q^h(x_0, u_0)$:

$$\sum_{j=0}^{K-1} \gamma^j r_{j+1}$$

- Furthermore, sample of $Q^h(x_k, u_k)$ at each step k :

$$Q^h(x_k, u_k) = \sum_{j=k}^{K-1} \gamma^{j-k} r_{j+1}$$

Monte Carlo policy evaluation (cont'd)



- To learn the expected value, run N trajectories (often called roll-outs)
- Estimated Q-value = **average** of the returns, e.g.

$$Q^h(x_0, u_0) = \frac{1}{N} \sum_{i=1}^N \sum_{j=0}^{K_i-1} \gamma^j r_{i,j+1}$$

Monte Carlo policy iteration

Monte Carlo policy iteration

```
for each iteration  $\ell$  do  
  run  $N$  trials applying  $h_\ell$   
  reset accumulator  $A(x, u)$ , counter  $C(x, u)$  to 0  
  for each step  $k$  of each trial  $i$  do  
     $A(x_k, u_k) \leftarrow A(x_k, u_k) + \sum_{j=k}^{K_i-1} \gamma^{j-k} r_{i,j+1}$  (return)  
     $C(x_k, u_k) \leftarrow C(x_k, u_k) + 1$   
  end for  
   $Q^{h_\ell}(x, u) \leftarrow A(x, u) / C(x, u)$   
   $h_{\ell+1}(x) \leftarrow \arg \max_u Q^{h_\ell}(x, u)$   
end for
```

Need for exploration

$$Q^h(x, u) \leftarrow A(x, u) / \mathbf{C(x, u)}$$

How to ensure $C(x, u) > 0$ – **information** about each (x, u) ?

- 1 Select representative **initial states** x_0
- 2 **Actions:**
 u_0 representative, sometimes different from $h(x_0)$
and in addition, perhaps:
 u_k representative, sometime different from $h(x_k)$



Exploration-exploitation

- **Exploration** needed:
actions different from the current policy
- **Exploitation** of current knowledge also needed:
current policy must be applied

Exploration-exploitation dilemma
– essential in all RL algorithms

(not just in MC)



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ϵ -greedy strategy

- Simple solution to the exploration-exploitation dilemma:

ϵ -greedy

$$u_k = \begin{cases} h(x_k) = \arg \max_u Q(x_k, u) & \text{with probability } (1 - \epsilon_k) \\ \text{a uniformly random action} & \text{w.p. } \epsilon_k \end{cases}$$

- Exploration probability $\epsilon_k \in (0, 1)$ usually decreased over time
- Main disadvantage: when exploring, actions are fully random, leading to poor performance

Softmax strategy

- Action selection:

$$u_k = u \text{ w.p. } \frac{e^{Q(x_k, u)/\tau_k}}{\sum_{u'} e^{Q(x_k, u')/\tau_k}}$$

where $\tau_k > 0$ is the **exploration temperature**

- Taking $\tau \rightarrow 0$, greedy selection recovered;
 $\tau \rightarrow \infty$ gives uniform random
- Compared to ε -greedy, better actions are more likely to be applied even when exploring

Bandit-based exploration



At single state, exploration modeled as **multi-armed bandit**:

- Action j = arm with reward distribution ρ_j , expectation μ_j
- Best arm (optimal action) has expected value μ^*
- At step k , we pull arm (try action) j_k , getting $r_k \sim \rho_{j_k}$
- **Objective:** After n pulls, small regret: $\sum_{k=1}^n \mu^* - \mu_{j_k}$

UCB algorithm

Popular algorithm: after n steps, pick arm with largest **upper confidence bound**:

$$b(j) = \hat{\mu}_j + \sqrt{\frac{3 \log n}{2n_j}}$$

where:

- $\hat{\mu}_j$ = mean of rewards observed for arm j so far
- n_j = how many times arm j was pulled

These are only a few simple methods, many others exist, e.g. Bayesian exploration, intrinsic rewards etc.



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DP perspective

- 1 Start from policy evaluation:

$$Q_{\tau+1}(x, u) \leftarrow \sum_{x'} f(x, u, x') [\rho(x, u, x') + \gamma Q_{\tau}(x', h(x'))]$$

- 2 Instead of model, use **transition sample** at each step k ,
 $(x_k, u_k, x_{k+1}, r_{k+1}, u_{k+1})$:

$$Q(x_k, u_k) \leftarrow r_{k+1} + \gamma Q(x_{k+1}, u_{k+1})$$

Note:

$$x_{k+1} \sim f(x_k, u_k, \cdot), r_{k+1} = \rho(x_k, u_k, x_{k+1}), u_{k+1} \sim h(x_{k+1})$$

- 3 Turn into **incremental update**:

$$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$$

$$[r_{k+1} + \gamma Q(x_{k+1}, u_{k+1}) - Q(x_k, u_k)]$$

$\alpha_k \in (0, 1]$ learning rate

Intermediate algorithm

Temporal differences for policy evaluation

for each trial **do**

 init x_0 , choose initial action u_0

repeat at each step k

 apply u_k , measure x_{k+1} , receive r_{k+1}

 choose **next** action $u_{k+1} \sim h(x_{k+1})$

$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$

$[r_{k+1} + \gamma Q(x_{k+1}, u_{k+1}) - Q(x_k, u_k)]$

until trial finished

end for

MC perspective

Temporal differences for policy h evaluation

for each trial **do**

...

repeat each step k

 apply u_k , measure x_{k+1} , receive r_{k+1}

$Q(x_k, u_k) \leftarrow \dots Q \dots$

until trial finished

end for

Monte Carlo

for each trial **do**

 execute trial

...

$Q(x, u) \leftarrow A(x, u) / C(x, u)$

end for



MC and DP perspectives

- Learn from online interaction: like MC, unlike DP
- Update after each transition, using previous Q-values: like DP, unlike MC

trial-based
transition-based

value iteration,
policy iteration

model-based

**temporal
differences**

Monte Carlo

model-free

Exploration-exploitation

choose next action $u_{k+1} \sim h(x_{k+1})$

- Information about $(x, u) \neq (x, h(x))$ needed
⇒ **exploration**
- h must be followed
⇒ **exploitation**

Policy improvement

- Previous algorithm: h fixed
- Improving h : simplest, after each transition, called optimistic policy improvement
- h implicit, greedy in Q
(update $Q \Rightarrow$ implicitly improve h)
- E.g. ε -greedy:

$$u_{k+1} = \begin{cases} \arg \max_u Q(x_{k+1}, u) & \text{w.p. } (1 - \varepsilon_{k+1}) \\ \text{uniformly random} & \text{w.p. } \varepsilon_{k+1} \end{cases}$$

SARSA

SARSA

for each trial **do**

init x_0

choose u_0 with exploration based on Q

repeat at each step k

apply u_k , measure x_{k+1} , receive r_{k+1}

choose u_{k+1} with exploration based on Q

$$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$$

$$[r_{k+1} + \gamma Q(x_{k+1}, u_{k+1}) - Q(x_k, u_k)]$$

until trial finished

end for

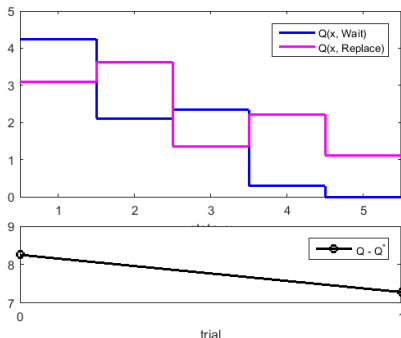
Origin of the name: $(x_k, u_k, r_{k+1}, x_{k+1}, u_{k+1}) =$
 (**S**tate, **A**ction, **R**eward, **S**tate, **A**ction)



Machine replacement: SARSA demo

Parameters: $\alpha = 0.1$, $\varepsilon = 0.3$ (constant), single trial
 $x_0 = 1$

SARSA, trial 1 completed



Q-learning

- 1 Similarly to SARSA, start from Q-iteration:

$$Q_{\ell+1}(x, u) \leftarrow \sum_{x'} f(x, u, x') [\rho(x, u, x') + \gamma \max_{u'} Q_{\ell}(x', u')]$$

- 2 Instead of model, use at each step k **transition sample**

$(x_k, u_k, x_{k+1}, r_{k+1})$:

$$Q(x_k, u_k) \leftarrow r_{k+1} + \gamma \max_{u'} Q(x_{k+1}, u')$$

Note: $x_{k+1} \sim f(x_k, u_k, \cdot)$, $r_{k+1} = \rho(x_k, u_k, x_{k+1})$

- 3 Turn into **incremental** update:

$$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$$

$$[r_{k+1} + \gamma \max_{u'} Q(x_{k+1}, u') - Q(x_k, u_k)]$$

Q-learning

Q-learning

for each trial **do**

 init x_0

repeat at each step k

choose u_k with exploration based on Q

 apply u_k , measure x_{k+1} , receive r_{k+1}

$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$

$[r_{k+1} + \gamma \max_{u'} Q(x_{k+1}, u') - Q(x_k, u_k)]$

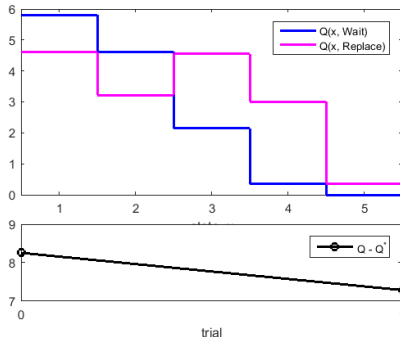
until trial finished

end for

Machine replacement: Q-learning demo

Parameters: $\alpha = 0.1$, $\varepsilon = 0.3$ (constant), single trial
 $x_0 = 1$

Q-learning, trial 1 completed



Convergence

Conditions for convergence to Q^*
in both SARSA and Q-learning:

- 1 All pairs (x, u) continue to be updated:
requires **exploration**, e.g. ε -greedy
- 2 Technical conditions on α_k (goes to 0, $\sum_{k=0}^{\infty} \alpha_k^2 = \text{finite}$,
but not too fast, $\sum_{k=0}^{\infty} \alpha_k \rightarrow \infty$)

In addition, for SARSA:

- 3 Policy must become greedy asymptotically
e.g. for ε -greedy, $\lim_{k \rightarrow \infty} \varepsilon_k = 0$



Discussion

SARSA **on-policy**

- Always updates towards the Q-function of the current policy

Q-learning **off-policy**

- Irrespective of the current policy, always updates towards optimal Q-function

Advantages of temporal differences

- Easy to understand and implement
- Low complexity \Rightarrow fast execution

Exploration and α_k sequence **greatly influence** performance

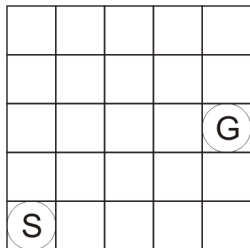


- 1 Introduction
- 2 Markov decision process & optimal solution
- 3 Dynamic programming, DP
- 4 Monte Carlo, MC
- 5 Exploration basics
- 6 Temporal differences, TD
- 7 Improving data efficiency

Motivation

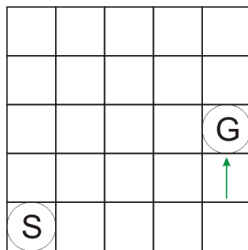
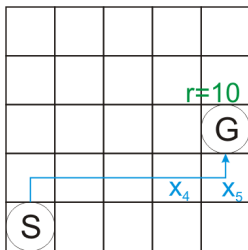
TD uses data inefficiently, and data is often costly.

Example:



- 2D gridworld navigation from **Start** to **Goal**
- Nonzero reward = 10 only on reaching G (terminal state)

Motivation (cont'd)



- Take SARSA with $\alpha = 1$; initialize $Q = 0$
- Updates along the trial on the left:

...

$$Q(x_4, u_4) = 0 + \gamma \cdot Q(x_5, u_5) = 0$$

$$Q(x_5, u_5) = 10 + \gamma \cdot 0 = 10$$

- A new transition from x_4 to x_5 (and hence a new trial) required to propagate the info to x_4 !

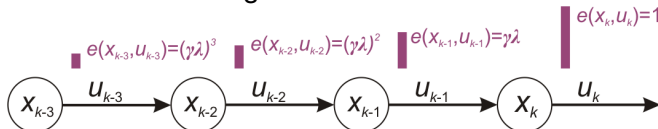
Ideas presented

- 1 Eligibility traces
- 2 Experience replay



Eligibility traces

- Idea: Leave a **trace** along the trial:



- $\lambda \in [0, 1]$ decay rate

Replacing traces

$e(x, u) \leftarrow 0$ for all x, u

for each step k do

$e(x, u) \leftarrow \lambda\gamma e(x, u)$ for all x, u

$e(x_k, u_k) \leftarrow 1$

end for

Example algorithm: SARSA(λ)

- Recall original SARSA only updates $Q(x_k, u_k)$:

$$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$$

$$[r_{k+1} + \gamma Q(x_{k+1}, u_{k+1}) - Q(x_k, u_k)]$$

- SARSA(λ) updates **all eligible pairs**:

$$Q(x, u) \leftarrow Q(x, u) + \alpha_k \cdot e(x, u) \cdot$$

$$[r_{k+1} + \gamma Q(x_{k+1}, u_{k+1}) - Q(x_k, u_k)] \quad \forall x, u$$



SARSA(λ)

SARSA(λ)

for each trial **do**

init x_0

$e(x, u) \leftarrow 0 \quad \forall x, u$

choose u_0 with exploration based on Q

repeat at each step k

apply u_k , measure x_{k+1} , receive r_{k+1}

choose u_{k+1} with exploration based on Q

$e(x, u) \leftarrow \lambda \gamma e(x, u) \quad \forall x, u$

$e(x_k, u_k) \leftarrow 1$

$Q(x, u) \leftarrow Q(x, u) + \alpha_k \cdot e(x, u) \cdot$

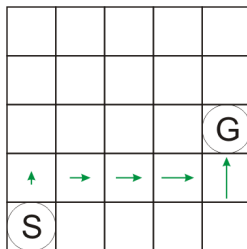
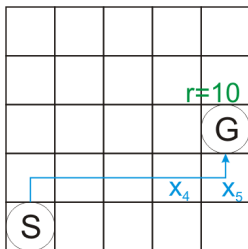
$[r_{k+1} + \gamma Q(x_{k+1}, u_{k+1}) - Q(x_k, u_k)]$ for all x, u

until trial finished

end for



Example: Effect of eligibility traces



- $\lambda = 0.7$
- Updates until x_4 : Q remains 0
- At x_5 , the entire trial gets updated:

$$Q(x_5, u_5) = 10 + \gamma 0 = 10$$

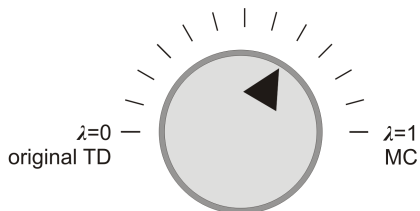
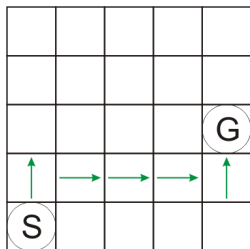
$$Q(x_4, u_4) = (\gamma \lambda)[10 + \gamma 0] = 3.5$$

$$Q(x_3, u_3) = (\gamma \lambda)^2[10 + \gamma 0] = 1.225$$

...

TD versus MC

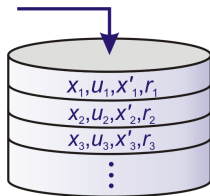
- $\lambda = 0 \Rightarrow$ recovers original algorithms, e.g. SARSA(0)
- $\lambda = 1 \Rightarrow$ TD becomes like MC



Typical values of λ are around 0.5 to 0.8

Experience replay

- Store each transition $(x_k, u_k, x_{k+1}, r_{k+1})$ in a database



- At each step, **replay** N transitions from the database (in addition to regular updates)

Q-learning with experience replay

Q-learning with experience replay

for each trial **do**

init x_0

repeat at each step k

apply u_k , measure x_{k+1} , receive r_{k+1}

$Q(x_k, u_k) \leftarrow Q(x_k, u_k) + \alpha_k \cdot$

$[r_{k+1} + \gamma \max_{u'} Q(x_{k+1}, u') - Q(x_k, u_k)]$

add $(x_k, u_k, x_{k+1}, r_{k+1})$ to database

ReplayExperience

until trial finished

end for

ReplayExperience procedure

ReplayExperience

loop N times

retrieve a transition (x, u, x', r) from database

$$Q(x, u) \leftarrow Q(x, u) + \alpha \cdot$$

$$[r + \gamma \max_{u'} Q(x', u') - Q(x, u)]$$

end loop

Retrieval order:

- 1 Backwards along trials, best for classical algos
- 2 Randomly, helps e.g. in deep RL
- 3 etc.



Textbooks

- Sutton & Barto, *Reinforcement Learning: An Introduction*, 1998 (+ ongoing 2nd edition, 2017).
- Bertsekas, *Dynamic Programming and Optimal Control*, vol. 2, 4th ed., 2012.
- Szepesvári, *Algorithms for Reinforcement Learning*, 2010.
- Buşoniu, Babuška, De Schutter, & Ernst, *Reinforcement Learning and Dynamic Programming Using Function Approximators*, 2010.

