Function Approximation, Deep Q Network, Rainbow

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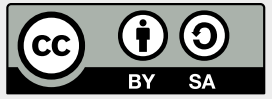


Function Approximation

We will approximate value function v and/or action-value function q, selecting it from a family of functions parametrized by a weight vector $\boldsymbol{w} \in \mathbb{R}^d$.

We denote the approximations as

$$\hat{v}(s; \boldsymbol{w}),$$
$$\hat{q}(s, a; \boldsymbol{w}).$$

We utilize the *Mean Squared Value Error* objective, denoted $\overline{VE}$:

$$\overline{VE}(\boldsymbol{w}) \stackrel{\text{def}}{=} \sum_{s \in \mathcal{S}} \mu(s) \left[v_\pi(s) - \hat{v}(s, \boldsymbol{w})\right]^2,$$

where the state distribution $\mu(s)$ is usually on-policy distribution.

Gradient and Semi-Gradient Methods

The functional approximation (i.e., the weight vector $\boldsymbol{w}$) is usually optimized using gradient methods, for example as

$$\begin{aligned}\boldsymbol{w}_{t+1} &\leftarrow \boldsymbol{w}_t - \tfrac{1}{2}\alpha\nabla_{\boldsymbol{w}_t}\big(v_\pi(S_t) - \hat{v}(S_t;\boldsymbol{w}_t)\big)^2 \\ &\leftarrow \boldsymbol{w}_t + \alpha\big(v_\pi(S_t) - \hat{v}(S_t;\boldsymbol{w}_t)\big)\nabla_{\boldsymbol{w}_t}\hat{v}(S_t;\boldsymbol{w}_t).\end{aligned}$$

As usual, the $v_\pi(S_t)$ is estimated by a suitable sample of a return:

- in Monte Carlo methods, we use episodic return G_t,
- in temporal difference methods, we employ bootstrapping and use one-step return

$$R_{t+1} + [\neg\text{done}] \cdot \gamma\hat{v}(S_{t+1};\boldsymbol{w})$$

  or an n-step return.

Temporal Difference Semi-Gradient Policy Evaluation

In TD methods, we again bootstrap the estimate $v_\pi(S_t)$ as $R_{t+1} + [\neg\text{done}] \cdot \gamma \hat{v}(S_{t+1}; \boldsymbol{w})$.

Semi-gradient TD(0) for estimating $\hat{v} \approx v_\pi$

Input: the policy π to be evaluated
Input: a differentiable function $\hat{v} : \mathcal{S}^+ \times \mathbb{R}^d \to \mathbb{R}$ 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'
    $\mathbf{w} \leftarrow \mathbf{w} + \alpha\big[R + \gamma\hat{v}(S',\mathbf{w}) - \hat{v}(S,\mathbf{w})\big]\nabla\hat{v}(S,\mathbf{w})$
    $S \leftarrow S'$
  until S is terminal

Algorithm 9.3 of "Reinforcement Learning: An Introduction, Second Edition".

Why Semi-Gradient TD

Note that the above algorithm is called **semi-gradient**, because it does not backpropagate through $\hat{v}(S_{t+1}; \boldsymbol{w})$:

$$\boldsymbol{w} \leftarrow \boldsymbol{w} + \alpha\Big(R_{t+1} + [\neg\text{done}] \cdot \gamma\hat{v}(S_{t+1}; \boldsymbol{w}) - \hat{v}(S_t; \boldsymbol{w})\Big)\nabla_{\boldsymbol{w}}\hat{v}(S_t; \boldsymbol{w}).$$

In other words, the above rule is in fact not an SGD update, because there does not exist a function $J(\boldsymbol{w})$, for which we would get the above update.

To sketch a proof, consider a linear $\hat{v}(S_t; \boldsymbol{w}) = \sum_i x(S_t)_i w_i$ and assume such a $J(\boldsymbol{w})$ exists. Then

$$\frac{\partial}{\partial w_i} J(\boldsymbol{w}) = \Big(R_{t+1} + \gamma\hat{v}(S_{t+1}; \boldsymbol{w}) - \hat{v}(S_t; \boldsymbol{w})\Big)x(S_t)_i.$$

Now considering second derivatives, we see they are not equal, which is a contradiction:

$$\frac{\partial}{\partial w_i}\frac{\partial}{\partial w_j} J(\boldsymbol{w}) = \Big(\gamma x(S_{t+1})_i - x(S_t)_i\Big)x(S_t)_j = \gamma x(S_{t+1})_i x(S_t)_j - x(S_t)_i x(S_t)_j$$
$$\frac{\partial}{\partial w_j}\frac{\partial}{\partial w_i} J(\boldsymbol{w}) = \Big(\gamma x(S_{t+1})_j - x(S_t)_j\Big)x(S_t)_i = \gamma x(S_{t+1})_j x(S_t)_i - x(S_t)_i x(S_t)_j$$

Temporal Difference Semi-Gradient Convergence

It can be proven (by using separate theory than for SGD) that the linear semi-gradient TD methods do converge.

However, they do not converge to the optimum of $\overline{VE}$. Instead, they converge to a different **TD fixed point** $\boldsymbol{w}_{\mathrm{TD}}$.

It can be proven that

$$\overline{VE}(\boldsymbol{w}_{\mathrm{TD}}) \leq \frac{1}{1-\gamma} \min_{\boldsymbol{w}} \overline{VE}(\boldsymbol{w}).$$

However, when γ is close to one, the multiplication factor in the above bound is quite large.

Temporal Difference Semi-Gradient Policy Evaluation

As before, we can utilize n-step TD methods.

n-step semi-gradient TD for estimating $\hat{v} \approx v_\pi$

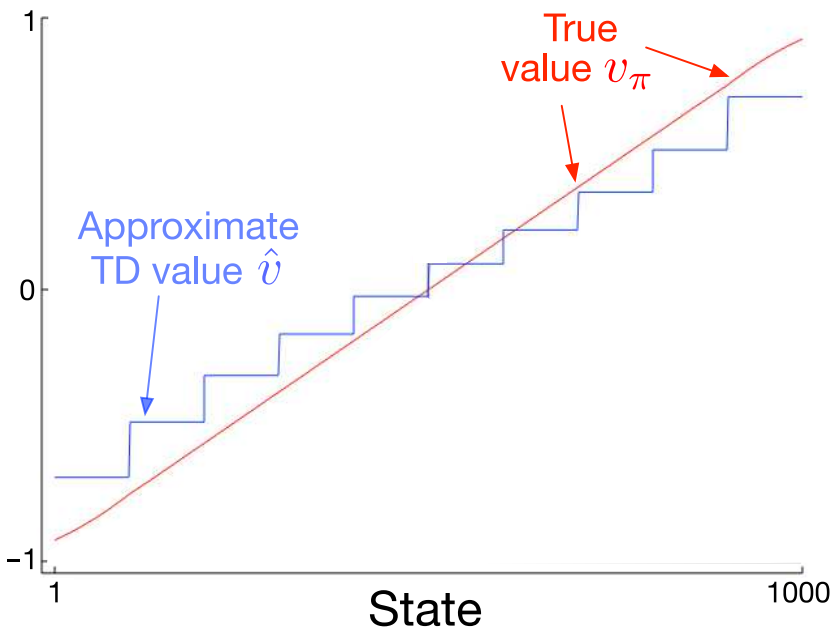
Input: the policy π to be evaluated
Input: a differentiable function $\hat{v} : \mathcal{S}^+ \times \mathbb{R}^d \to \mathbb{R}$ such that $\hat{v}(\text{terminal},\cdot) = 0$
Algorithm parameters: step size $\alpha > 0$, a positive integer n
Initialize value-function weights $\mathbf{w}$ arbitrarily (e.g., $\mathbf{w} = \mathbf{0}$)
All store and access operations (S_t and R_t) can take their index mod $n + 1$

Loop for each episode:
  Initialize and store $S_0 \neq$ terminal
  $T \leftarrow \infty$
  Loop for $t = 0, 1, 2, \ldots$:
  | If $t < T$, then:
  | Take an action according to $\pi(\cdot|S_t)$
  | Observe and store the next reward as R_{t+1} and the next state as S_{t+1}
  | If S_{t+1} is terminal, then $T \leftarrow t + 1$
  | $\tau \leftarrow t - n + 1$ (τ is the time whose state's estimate is being updated)
  | If $\tau \geq 0$:
  | $G \leftarrow \sum_{i=\tau+1}^{\min(\tau+n,T)} \gamma^{i-\tau-1} R_i$
  | If $\tau + n < T$, then: $G \leftarrow G + \gamma^n \hat{v}(S_{\tau+n},\mathbf{w})$ $(G_{\tau:\tau+n})$
  | $\mathbf{w} \leftarrow \mathbf{w} + \alpha \left[G - \hat{v}(S_\tau,\mathbf{w})\right] \nabla \hat{v}(S_\tau,\mathbf{w})$
  Until $\tau = T - 1$

Algorithm 9.5 of "Reinforcement Learning: An Introduction, Second Edition".

Temporal Difference Semi-Gradient Policy Evaluation

On the left, the results of one-step TD(0) algorithm are presented. The effect of increasing n in an n-step variant is displayed on the right.



Average RMS error over 1000 states and first 10 episodes

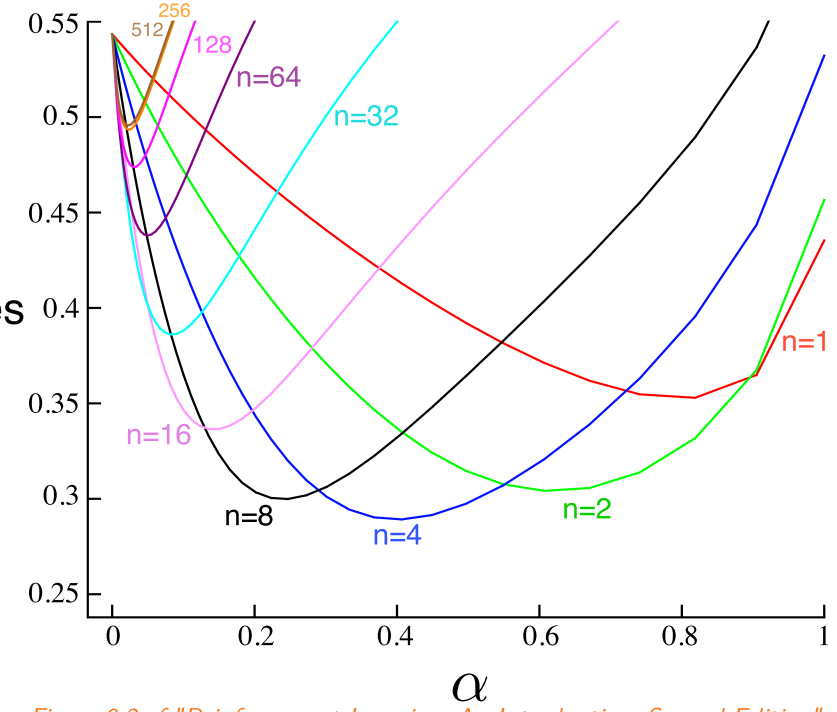


Figure 9.2 of "Reinforcement Learning: An Introduction, Second Edition".

Sarsa with Function Approximation

Until now, we talked only about policy evaluation. Naturally, we can extend it to a full Sarsa algorithm:

Episodic Semi-gradient Sarsa for Estimating $\hat{q} \approx q_*$

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

Loop for each episode:
  $S, A \leftarrow$ initial state and action of episode (e.g., ε-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)
    $\mathbf{w} \leftarrow \mathbf{w} + \alpha\big[R + \gamma\hat{q}(S', A', \mathbf{w}) - \hat{q}(S, A, \mathbf{w})\big]\nabla\hat{q}(S, A, \mathbf{w})$
    $S \leftarrow S'$
    $A \leftarrow A'$

Modified from Algorithm 10.1 of "Reinforcement Learning: An Introduction, Second Edition".

Sarsa with Function Approximation

Additionally, we can incorporate n-step returns:

Episodic semi-gradient n-step Sarsa for estimating $\hat{q} \approx q_*$ or q_π

Input: a differentiable action-value function parameterization $\hat{q} : \mathcal{S} \times \mathcal{A} \times \mathbb{R}^d \to \mathbb{R}$
Input: a policy π (if estimating q_π)
Algorithm parameters: step size $\alpha > 0$, small $\varepsilon > 0$, a positive integer n
Initialize action-value function weights $\mathbf{w} \in \mathbb{R}^d$ arbitrarily (e.g., $\mathbf{w} = \mathbf{0}$)
All store and access operations (S_t, A_t, and R_t) can take their index mod $n + 1$

Loop for each episode:
  Initialize and store $S_0 \neq$ terminal
  Select and store an action $A_0 \sim \pi(\cdot|S_0)$ or ε-greedy wrt $\hat{q}(S_0, \cdot, \mathbf{w})$
  $T \leftarrow \infty$
  Loop for $t = 0, 1, 2, \ldots$:
    If $t < T$, then:
      Take action A_t
      Observe and store the next reward as R_{t+1} and the next state as S_{t+1}
      If S_{t+1} is terminal, then:
        $T \leftarrow t + 1$
      else:
        Select and store $A_{t+1} \sim \pi(\cdot|S_{t+1})$ or ε-greedy wrt $\hat{q}(S_{t+1}, \cdot, \mathbf{w})$
    $\tau \leftarrow t - n + 1$ (τ is the time whose estimate is being updated)
    If $\tau \geq 0$:
      $G \leftarrow \sum_{i=\tau+1}^{\min(\tau+n,T)} \gamma^{i-\tau-1} R_i$
      If $\tau + n < T$, then $G \leftarrow G + \gamma^n \hat{q}(S_{\tau+n}, A_{\tau+n}, \mathbf{w})$ $(G_{\tau:\tau+n})$
      $\mathbf{w} \leftarrow \mathbf{w} + \alpha \left[G - \hat{q}(S_\tau, A_\tau, \mathbf{w})\right] \nabla \hat{q}(S_\tau, A_\tau, \mathbf{w})$
  Until $\tau = T - 1$

Modified from Algorithm 10.2 of "Reinforcement Learning: An Introduction, Second Edition".

Mountain Car Example

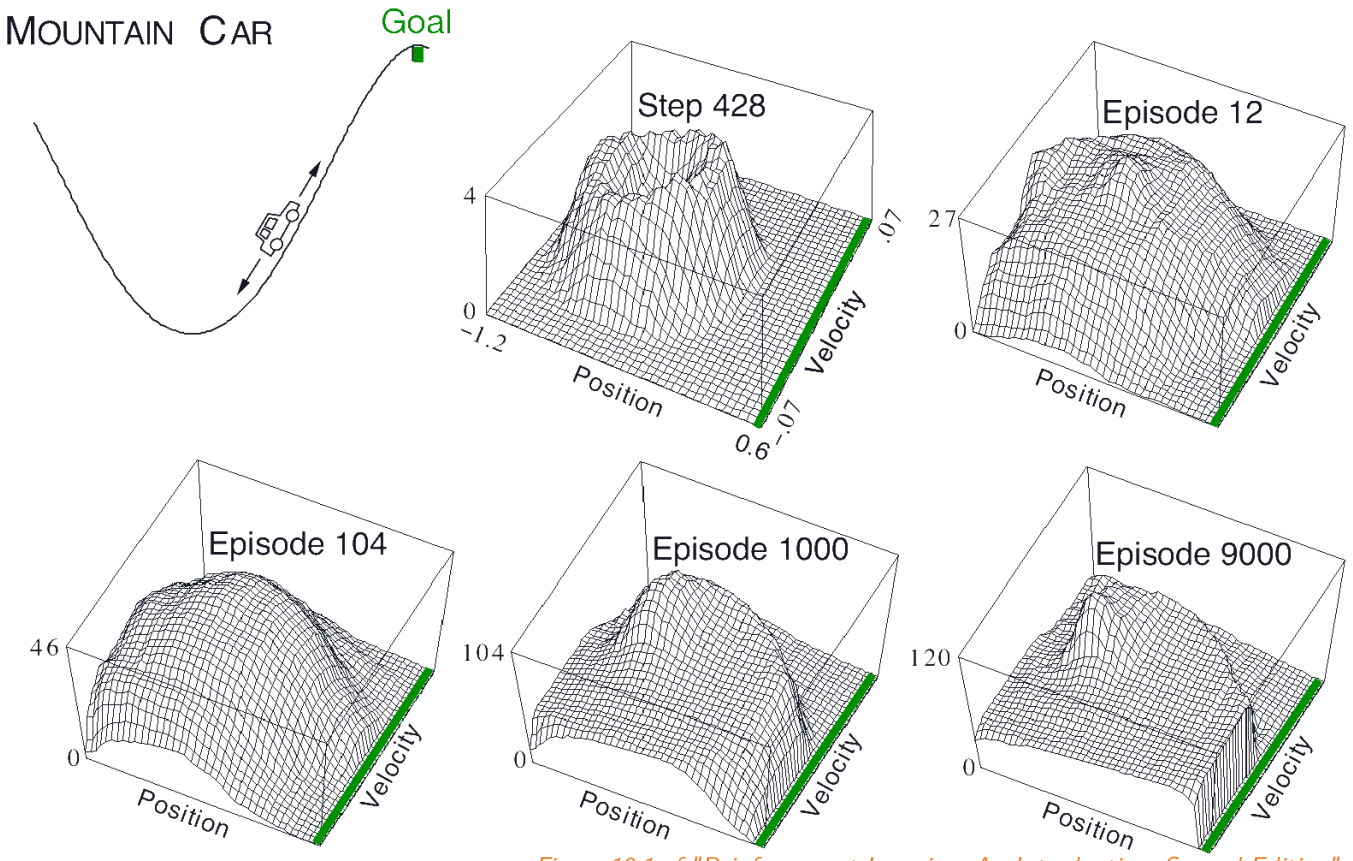


Figure 10.1 of "Reinforcement Learning: An Introduction, Second Edition".

The performances are for semi-gradient Sarsa(λ) algorithm (which we did not talked about yet) with tile coding of 8 overlapping tiles covering position and velocity, with offsets of $(1, 3)$.

Mountain Car Example

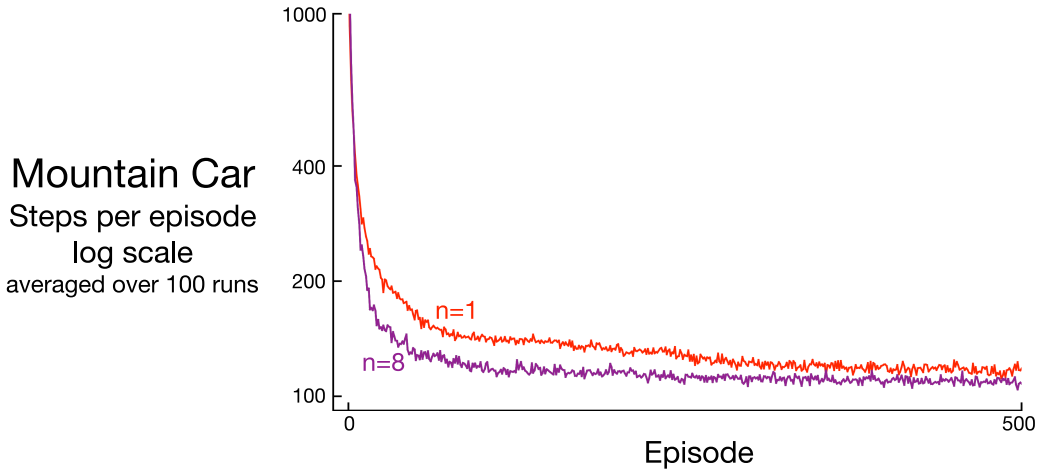


Figure 10.3 of "Reinforcement Learning: An Introduction, Second Edition".

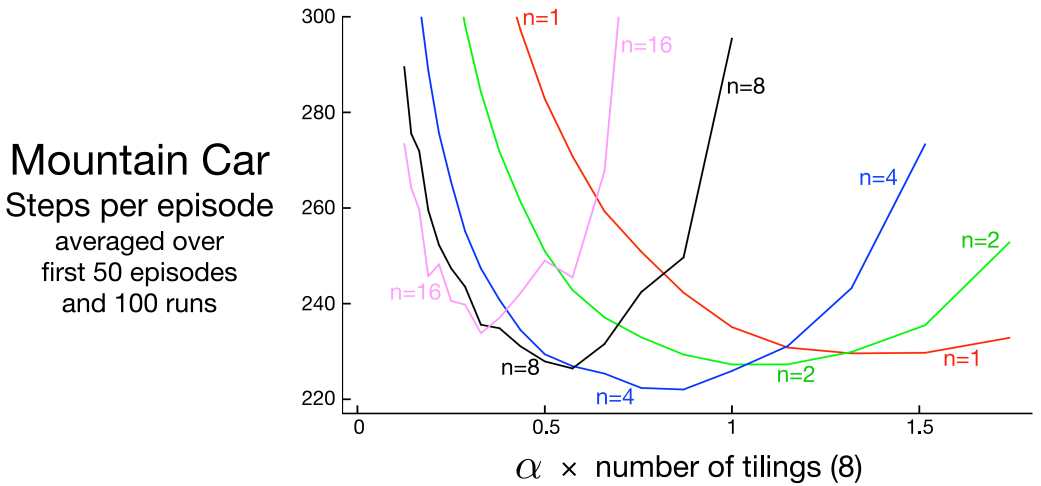


Figure 10.4 of "Reinforcement Learning: An Introduction, Second Edition".

Off-policy Divergence With Function Approximation

Consider a deterministic transition between two states whose values are computed using the same weight:

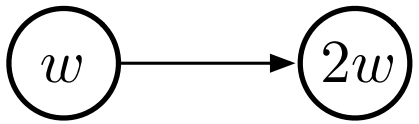


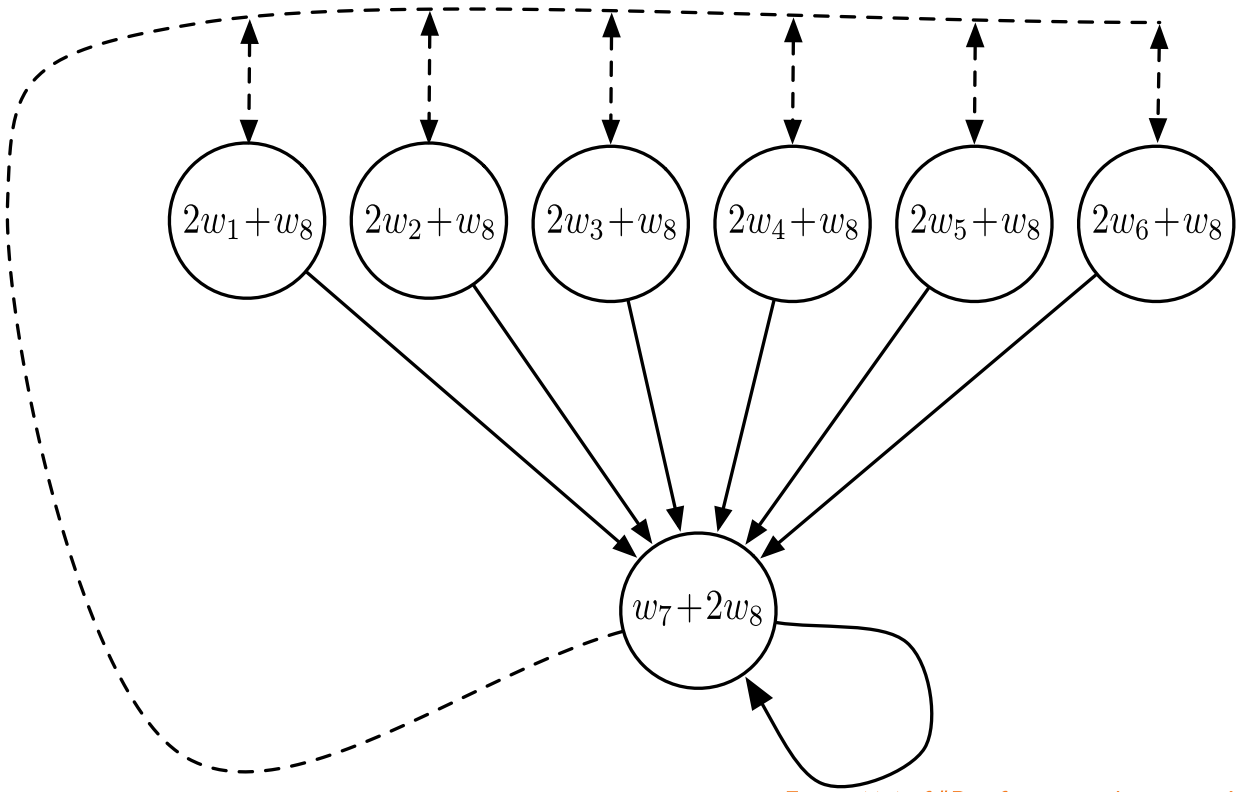
Figure from Section 11.2 of "Reinforcement Learning: An Introduction, Second Edition".

- If initially $w = 10$, the TD error will be also 10 (or nearly 10 if $\gamma < 1$).
- If for example $\alpha = 0.1$, w will be increased to 11 (by 10%).
- This process can continue indefinitely.

However, the problem arises only in off-policy setting, where we do not decrease value of the second state from further observation.

Off-policy Divergence With Function Approximation

The previous idea can be implemented for instance by the following **Baird's counterexample**:



$\pi(\text{solid}|\cdot) = 1$

$b(\text{dashed}|\cdot) = 6/7$

$b(\text{solid}|\cdot) = 1/7$

$\gamma = 0.99$

Figure 11.1 of "Reinforcement Learning: An Introduction, Second Edition".

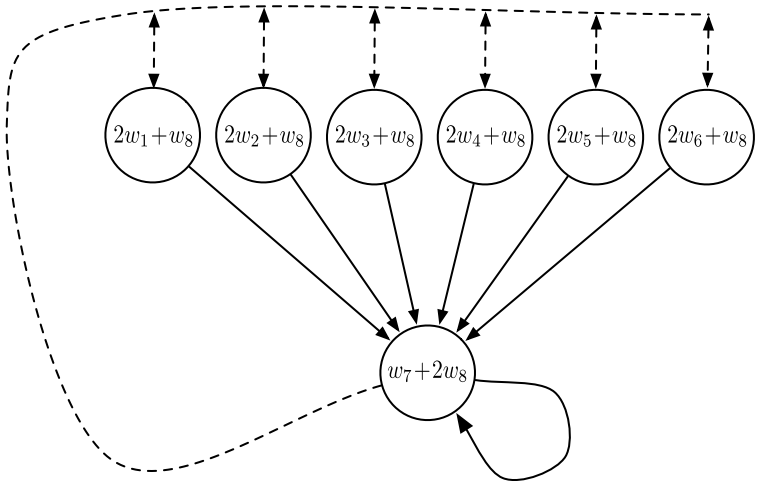
The rewards are zero everywhere, so the value function is also zero everywhere. We assume the initial values of weights are 1, except for $w_7 = 10$, and that the learning rate $\alpha = 0.01$.

Off-policy Divergence With Function Approximation

For off-policy semi-gradient Sarsa, or even for off-policy dynamic-programming update (where we compute expectation over all following states and actions), the weights diverge to $+\infty$.

Using on-policy distribution converges fine.

$$\boldsymbol{w} \leftarrow \boldsymbol{w} + \frac{\alpha}{|\mathcal{S}|} \sum_{\boldsymbol{s}} \Big(\mathbb{E}_\pi \left[R_{t+1} + \gamma \hat{v}(S_{t+1}; \boldsymbol{w}) | S_t = \boldsymbol{s} \right] - \hat{v}(\boldsymbol{s}; \boldsymbol{w}) \Big) \nabla \hat{v}(\boldsymbol{s}; \boldsymbol{w})$$



$\pi(\text{solid}|\cdot) = 1$

$b(\text{dashed}|\cdot) = 6/7$

$b(\text{solid}|\cdot) = 1/7$

$\gamma = 0.99$

Figure 11.1 of "Reinforcement Learning: An Introduction, Second Edition".

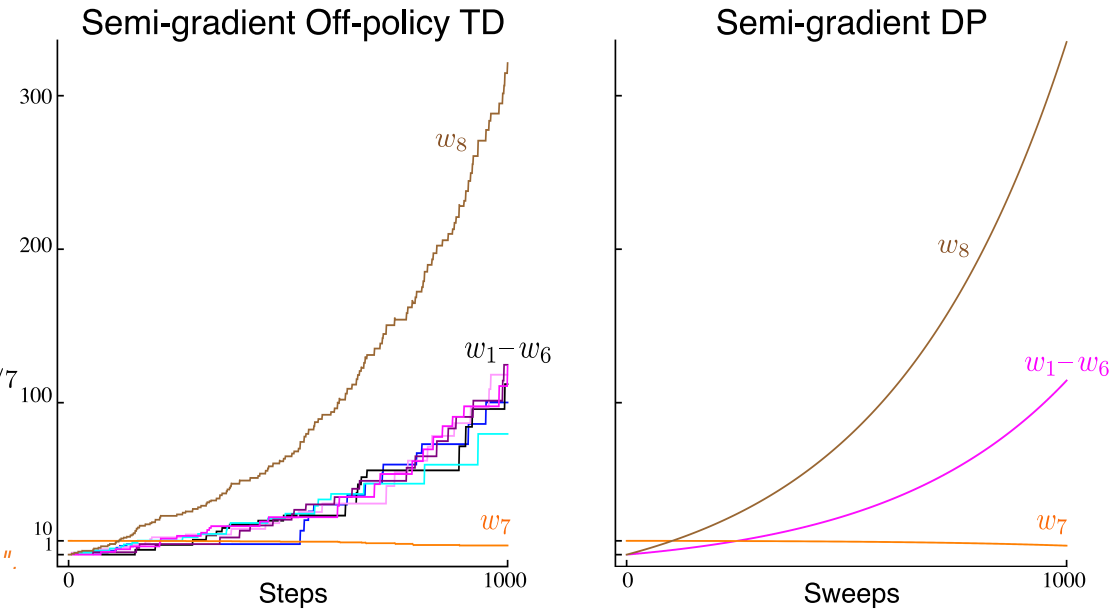


Figure 11.2 of "Reinforcement Learning: An Introduction, Second Edition".

Off-policy Divergence With Function Approximation

The divergence can happen when all following elements are combined:

- functional approximation;
- bootstrapping;
- off-policy training.

In the Sutton's and Barto's book, these are called **the deadly triad**.

Deep Q Networks

Volodymyr Mnih et al.: *Playing Atari with Deep Reinforcement Learning* (Dec 2013 on arXiv), in Feb 2015 accepted in Nature as *Human-level control through deep reinforcement learning*.

Off-policy Q-learning algorithm with a convolutional neural network function approximation of action-value function.

Training can be extremely brittle (and can even diverge as shown earlier).

Deep Q Network

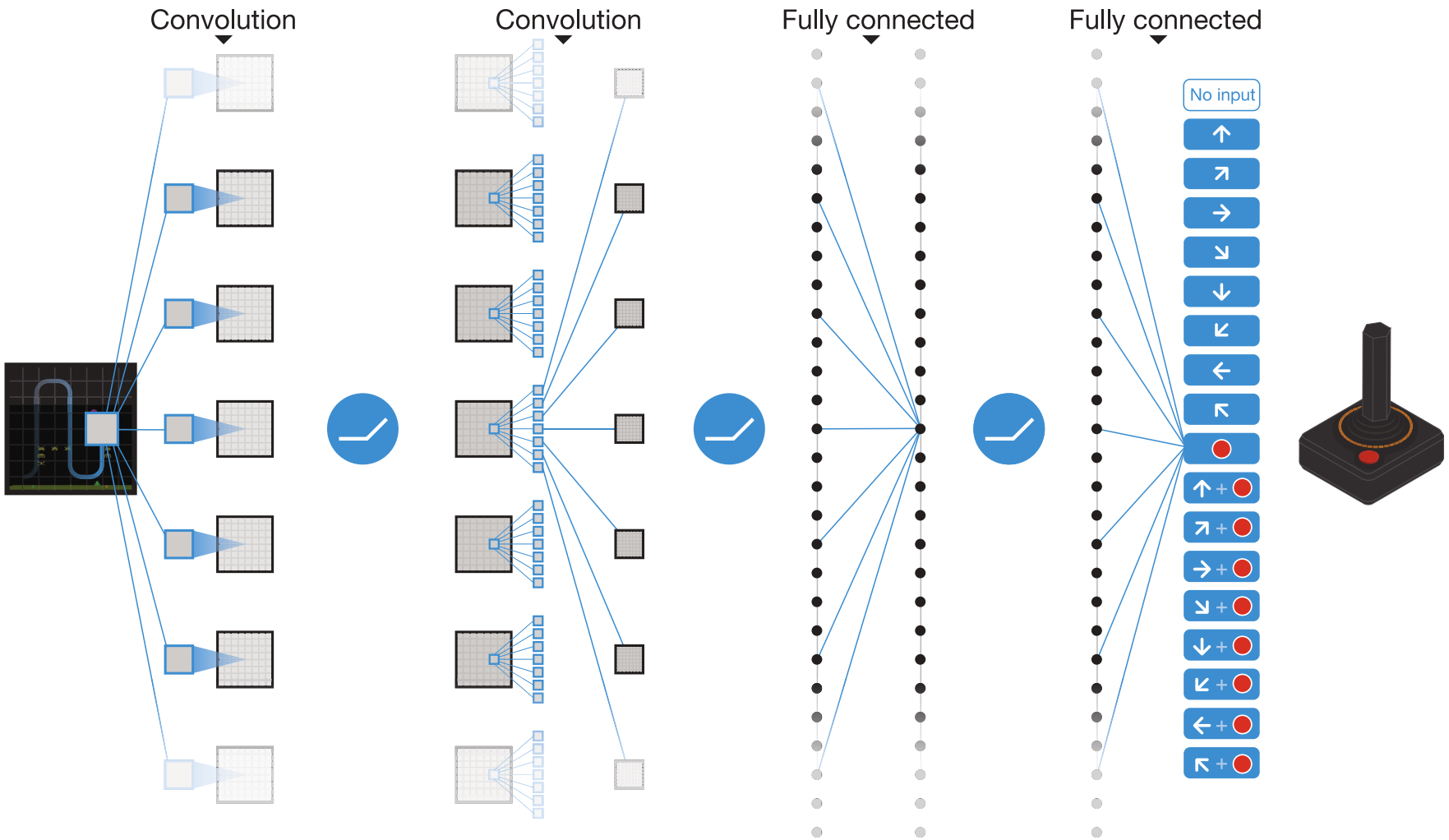


Figure 1 of "Human-level control through deep reinforcement learning" by Volodymyr Mnih et al.

Deep Q Networks

- Preprocessing: 210×160 128-color images are converted to grayscale and then resized to 84×84.
- **Frame skipping** technique is used, i.e., only every 4^{th} frame (out of 60 per second) is considered, and the selected action is repeated on the other frames.
- **Frame stacking** is utilizied – the input to the network are the last 4 frames (considering only the frames kept by frame skipping), i.e., the network inpus is an image with 4 channels.
- The network is fairly standard, performing
  - 32 filters of size 8×8 with stride 4 and ReLU,
  - 64 filters of size 4×4 with stride 2 and ReLU,
  - 64 filters of size 3×3 with stride 1 and ReLU,
  - fully connected layer with 512 units and ReLU,
  - output layer with 18 output units (one for each action)

- Network is trained with RMSProp to minimize the following loss:

$$\mathcal{L} \stackrel{\text{def}}{=} \mathbb{E}_{(s,a,r,s')\sim\text{data}}\left[\left(r + [\neg\text{done}]\cdot\gamma\max_{a'} Q(s',a';\bar{\boldsymbol{\theta}}) - Q(s,a;\boldsymbol{\theta})\right)^2\right].$$

- An ε-greedy behavior policy is utilized (starts at $\varepsilon = 1$ and gradually decreases to 0.1).

Important improvements:

- **experience replay**: the generated episodes are stored in a buffer as (s, a, r, s') quadruples, and for training a transition is sampled uniformly (off-policy training);
- separate **target network $\bar{\boldsymbol{\theta}}$**: to prevent instabilities, a separate *target network* is used to estimate one-step returns. The weights are not trained, but copied from the trained network after a fixed number of gradient updates;
- reward clipping: because rewards have wildly different scale in different games, all positive rewards are replaced by $+1$ and negative by -1; life loss is used as an end of episode.
  - furthermore, $\left(r + [\neg\text{done}]\cdot\gamma\max_{a'} Q(s',a';\bar{\boldsymbol{\theta}}) - Q(s,a;\boldsymbol{\theta})\right)$ is also clipped to $[-1,1]$ (i.e., a smooth_{L_1} loss or Huber loss).

Algorithm 1: deep Q-learning with experience replay.

Initialize replay memory D to capacity N

Initialize action-value function Q with random weights θ

Initialize target action-value function $\hat{Q}$ with weights $\theta^- = \theta$

For episode $= 1, M$ **do**

  Initialize sequence $s_1 = \{x_1\}$ and preprocessed sequence $\phi_1 = \phi(s_1)$

  For $t = 1$,T **do**

    With probability ε select a random action a_t

    otherwise select $a_t = \text{argmax}_a Q(\phi(s_t), a; \theta)$

    Execute action a_t in emulator and observe reward r_t and image x_{t+1}

    Set $s_{t+1} = s_t, a_t, x_{t+1}$ and preprocess $\phi_{t+1} = \phi(s_{t+1})$

    Store transition $\left(\phi_t, a_t, r_t, \phi_{t+1}\right)$ in D

    Sample random minibatch of transitions $\left(\phi_j, a_j, r_j, \phi_{j+1}\right)$ from D

    Set $y_j = \begin{cases} r_j & \text{if episode terminates at step j+1} \\ r_j + \gamma \max_{a'} \hat{Q}\left(\phi_{j+1}, a'; \theta^-\right) & \text{otherwise} \end{cases}$

    Perform a gradient descent step on $\left(y_j - Q\left(\phi_j, a_j; \theta\right)\right)^2$ with respect to the network parameters θ

    Every C steps reset $\hat{Q} = Q$

  End For

End For

Algorithm 1 of "Human-level control through deep reinforcement learning" by Volodymyr Mnih et al.

Deep Q Network

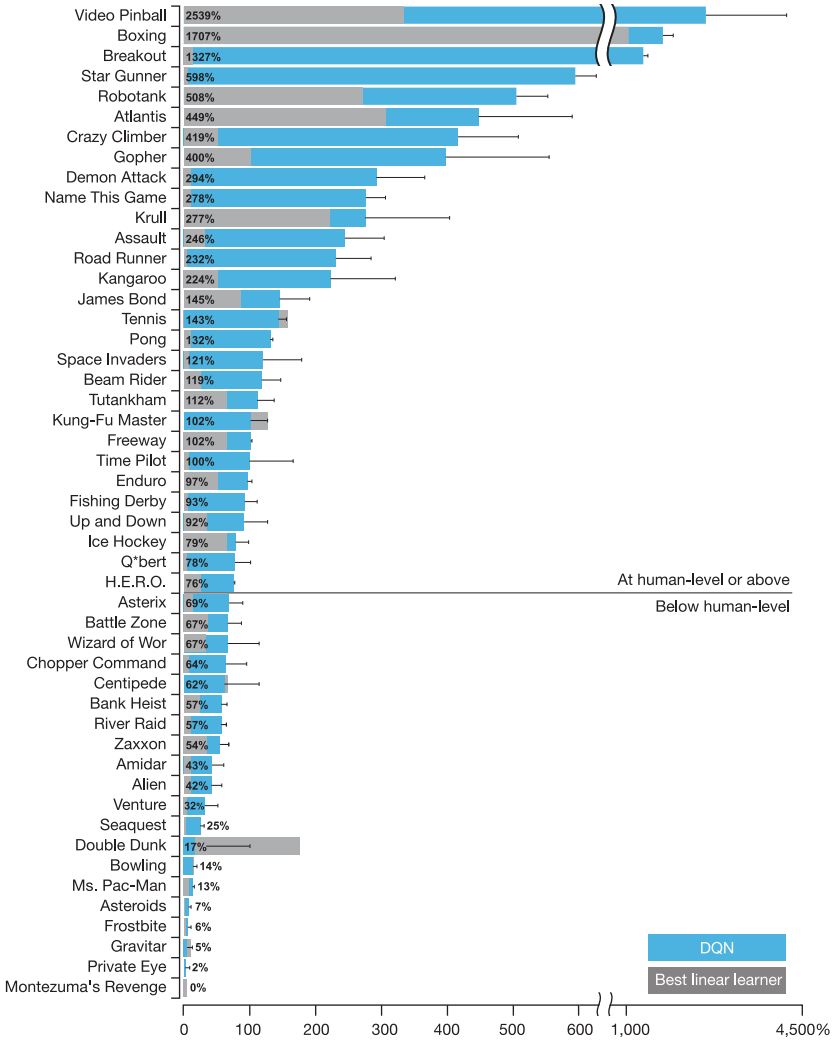
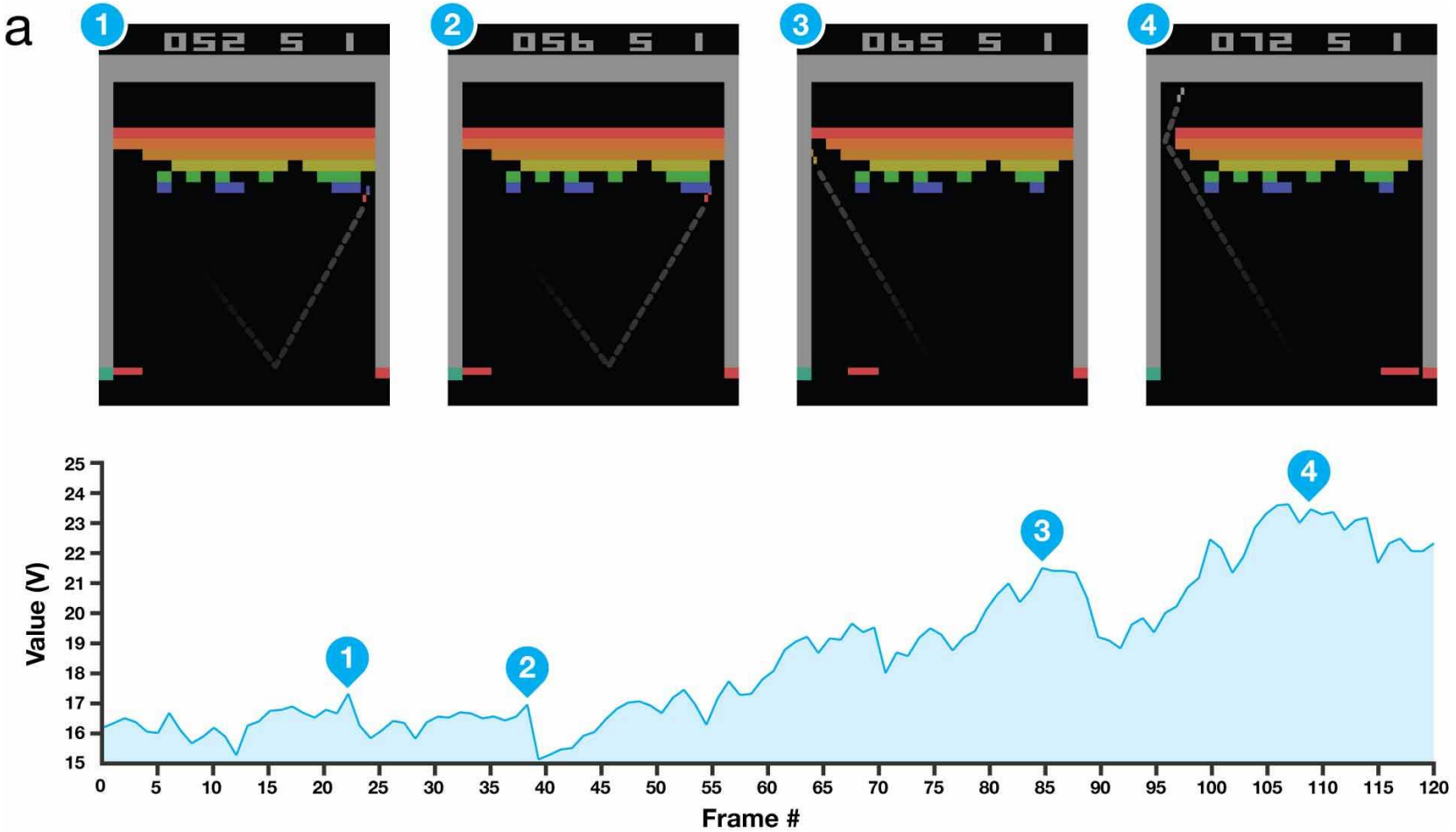


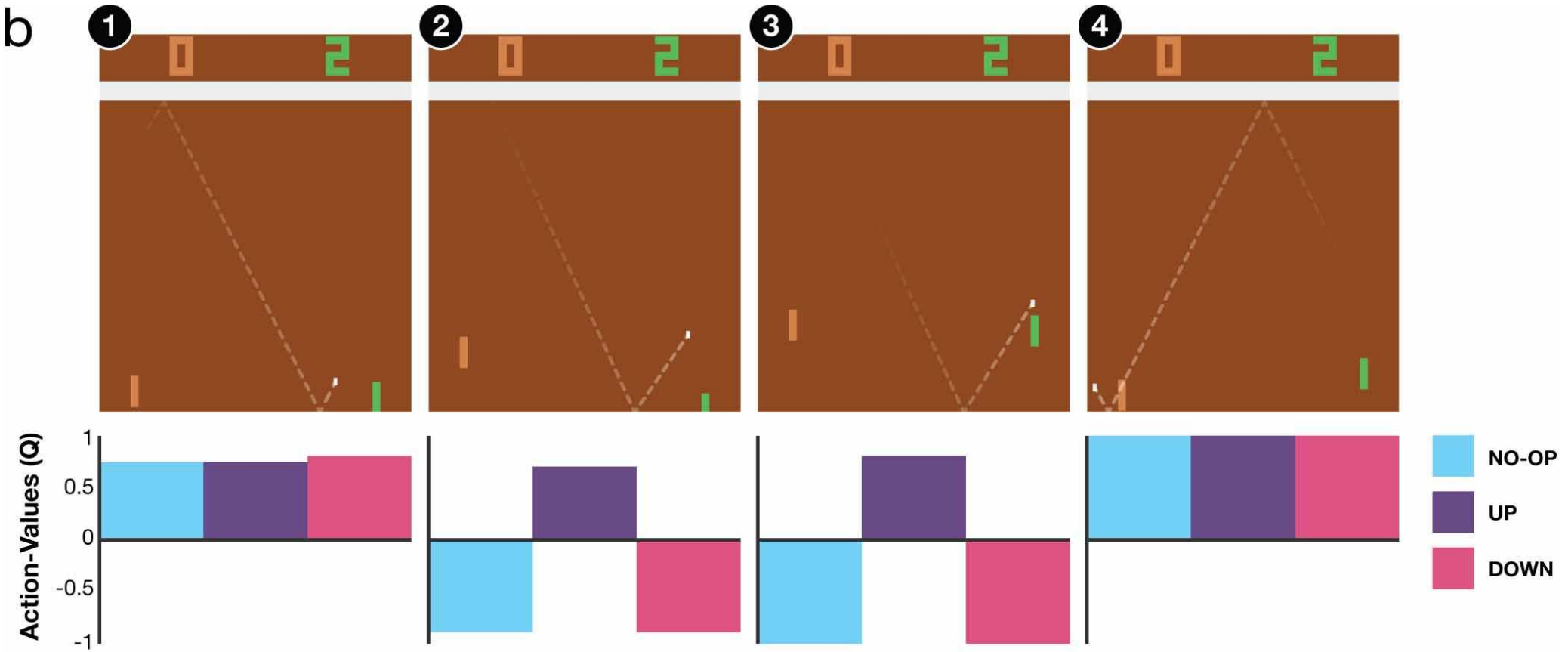
Figure 3 of "Human-level control through deep reinforcement learning" by Volodymyr Mnih et al.

Deep Q Network



Extended Data Figure 2a of "Human-level control through deep reinforcement learning" by Volodymyr Mnih et al.

Deep Q Network



Extended Data Figure 2b of "Human-level control through deep reinforcement learning" by Volodymyr Mnih et al.

Deep Q Networks Hyperparameters

Hyperparameter	Value
minibatch size	32
replay buffer size	1M
target network update frequency	10k
discount factor	0.99
training frames	50M
RMSProp learning rate and both momentums	0.00025, 0.95
initial ε, final ε (linear decay) and frame of final ε	1.0, 0.1, 1M
replay start size	50k
no-op max	30

Rainbow

There have been many suggested improvements to the DQN architecture. In the end of 2017, the *Rainbow: Combining Improvements in Deep Reinforcement Learning* paper combines 6 of them into a single architecture they call **Rainbow**.

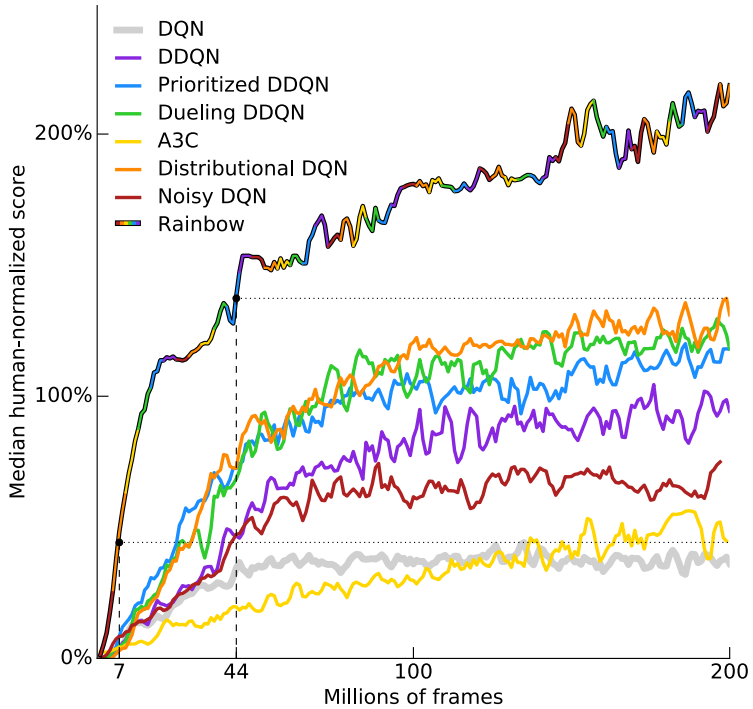


Figure 1 of "Rainbow: Combining Improvements in Deep Reinforcement Learning" by Matteo Hessel et al.

Q-learning and Maximization Bias

Because behaviour policy in Q-learning is ε-greedy variant of the target policy, the same samples (up to ε-greedy) determine both the maximizing action and estimate its value.

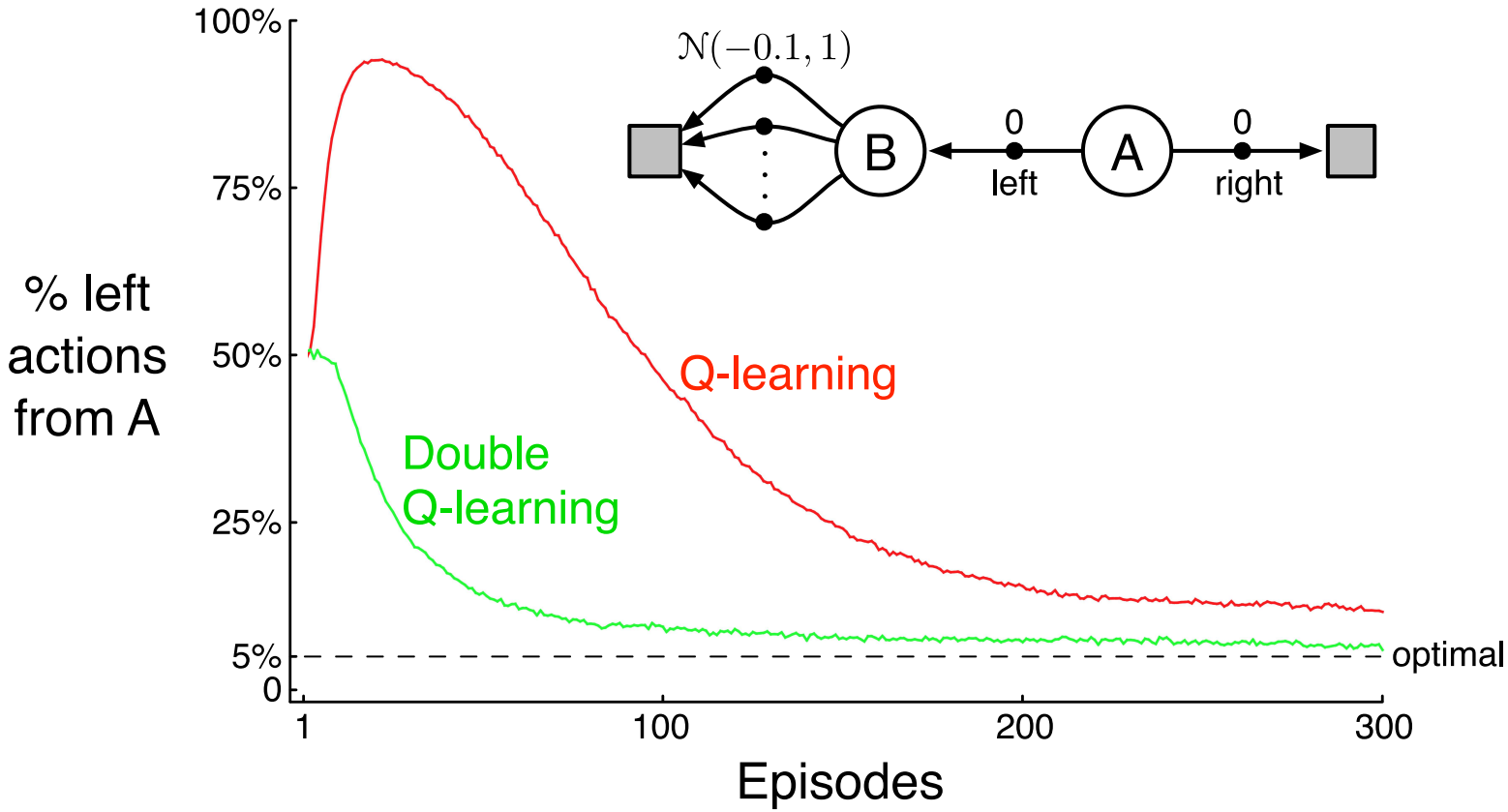


Figure 6.5 of "Reinforcement Learning: An Introduction, Second Edition".

Double Q-learning

Double Q-learning, for estimating $Q_1 \approx Q_2 \approx q_*$

Algorithm parameters: step size $\alpha \in (0, 1]$, small $\varepsilon > 0$
Initialize $Q_1(s, a)$ and $Q_2(s, a)$, for all $s \in \mathcal{S}, a \in \mathcal{A}(s)$, such that $Q(\mathit{terminal}, \cdot) = 0$

Loop for each episode:
  Initialize S
  Loop for each step of episode:
    Choose A from S using the policy ε-greedy in $Q_1 + Q_2$
    Take action A, observe R, S'
    With 0.5 probabilility:

$$Q_1(S, A) \leftarrow Q_1(S, A) + \alpha\Big(R + \gamma Q_2\big(S', \operatorname{argmax}_a Q_1(S', a)\big) - Q_1(S, A)\Big)$$

    else:

$$Q_2(S, A) \leftarrow Q_2(S, A) + \alpha\Big(R + \gamma Q_1\big(S', \operatorname{argmax}_a Q_2(S', a)\big) - Q_2(S, A)\Big)$$

    $S \leftarrow S'$
  until S is terminal

Modification of Algorithm 6.7 of "Reinforcement Learning: An Introduction, Second Edition" (replacing S+ by S).

Double Q-learning

Similarly to double Q-learning, instead of

$$r + \gamma \max_{a'} Q(s', a'; \bar{\boldsymbol{\theta}}) - Q(s, a; \boldsymbol{\theta}),$$

we minimize

$$r + \gamma Q(s', \arg\max_{a'} Q(s', a'; \boldsymbol{\theta}); \bar{\boldsymbol{\theta}}) - Q(s, a; \boldsymbol{\theta}).$$

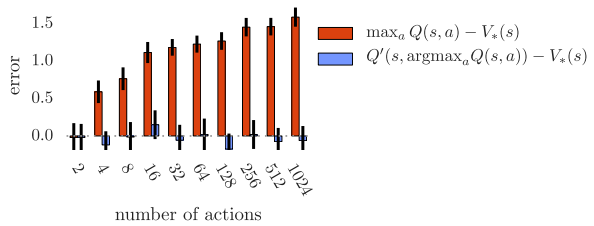


Figure 1: The orange bars show the bias in a single Q-learning update when the action values are $Q(s,a) = V_*(s) + \epsilon_a$ and the errors $\{\epsilon_a\}_{a=1}^m$ are independent standard normal random variables. The second set of action values Q', used for the blue bars, was generated identically and independently. All bars are the average of 100 repetitions.

Figure 1 of "Deep Reinforcement Learning with Double Q-learning" by Hado van Hasselt et al.

Double Q-learning

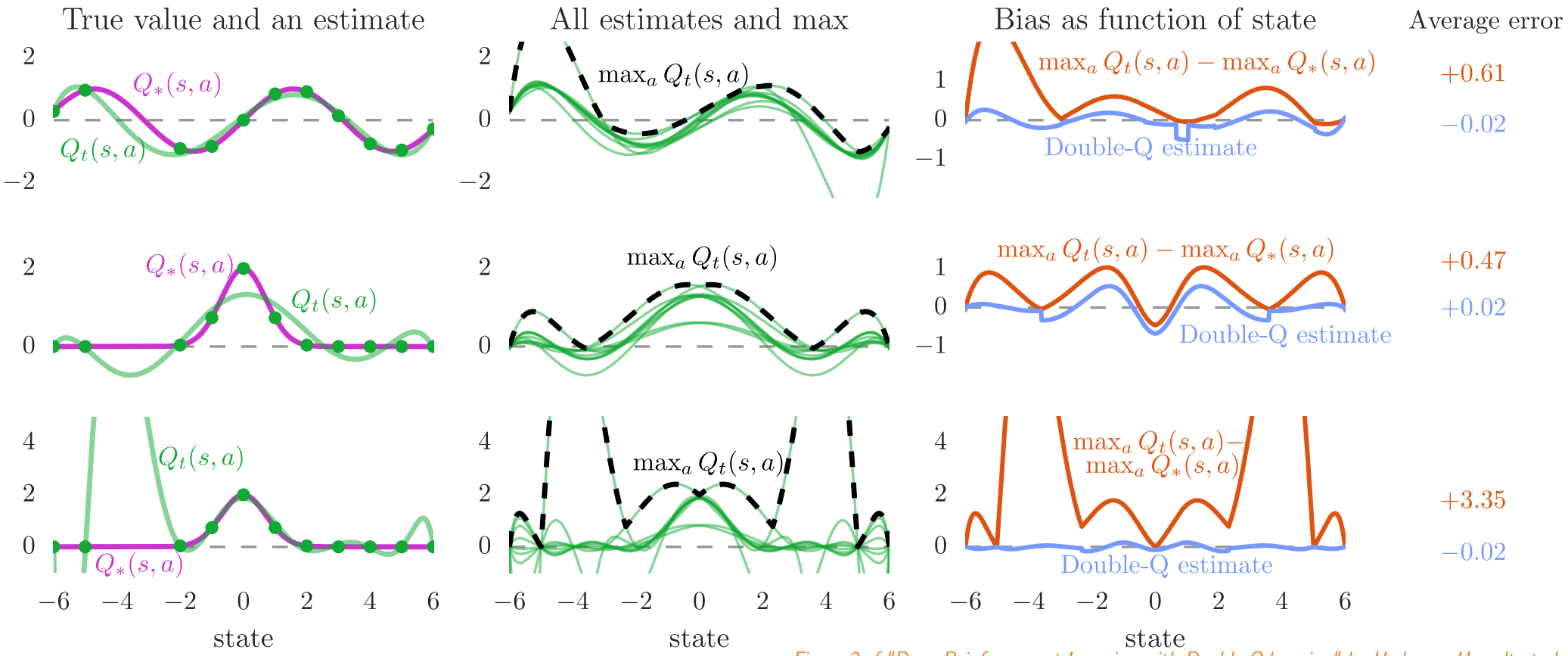


Figure 2 of "Deep Reinforcement Learning with Double Q-learning" by Hado van Hasselt et al.

Double Q-learning

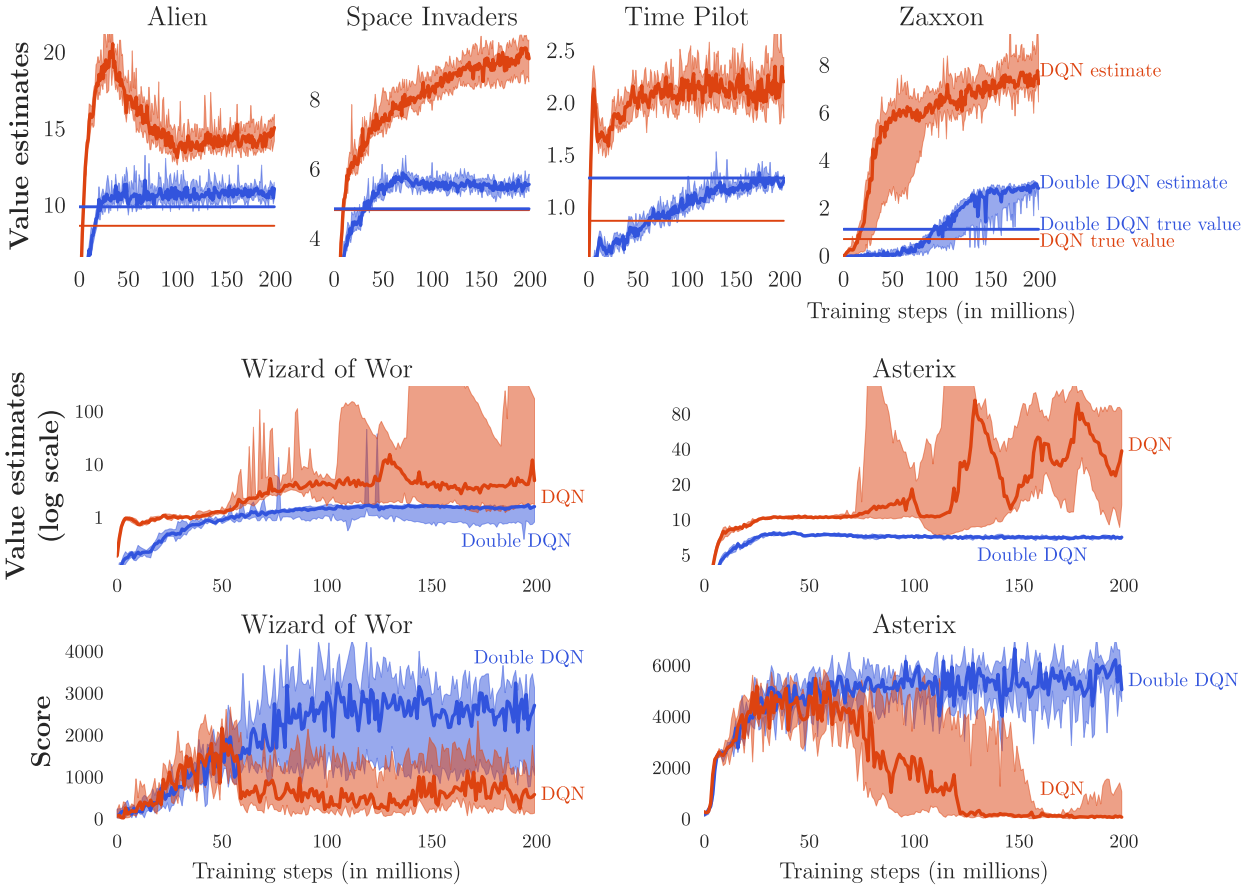


Figure 3 of "Deep Reinforcement Learning with Double Q-learning" by Hado van Hasselt et al.

Double Q-learning

Performance on episodes taking at most 5 minutes and no-op starts on 49 games:

	DQN	Double DQN
Median	93.5%	114.7%
Mean	241.1%	330.3%

Table 1 of "Deep Reinforcement Learning with Double Q-learning" by Hado van Hasselt et al.

Performance on episodes taking at most 30 minutes and using human starts on 49 games:

	DQN	Double DQN	Double DQN (tuned)
Median	47.5%	88.4%	116.7%
Mean	122.0%	273.1%	475.2%

Table 2 of "Deep Reinforcement Learning with Double Q-learning" by Hado van Hasselt et al.

Prioritized Replay

Instead of sampling the transitions uniformly from the replay buffer, we instead prefer those with a large TD error. Therefore, we sample transitions according to their probability

$$p_t \propto \left| r + \gamma \max_{a'} Q(s', a'; \bar{\boldsymbol{\theta}}) - Q(s, a; \boldsymbol{\theta}) \right|^{\omega},$$

where ω controls the shape of the distribution (which is uniform for $\omega = 0$ and corresponds to TD error for $\omega = 1$).

New transitions are inserted into the replay buffer with maximum probability to support exploration of all encountered transitions.

When combined with DDQN, the probabilities are naturally computed as

$$p_t \propto \left| r + \gamma Q(s', \arg\max_{a'} Q(s', a'; \boldsymbol{\theta}); \bar{\boldsymbol{\theta}}) - Q(s, a; \boldsymbol{\theta}) \right|^{\omega},$$

Prioritized Replay

Because we now sample transitions according to p_t instead of uniformly, on-policy distribution and sampling distribution differ. To compensate, we therefore utilize importance sampling with ratio

$$\rho_t = \left(\frac{1/N}{p_t}\right)^{\beta}.$$

The authors utilize in fact "for stability reasons"

$$\rho_t \Big/ \max_i \rho_i.$$

Prioritized Replay

Algorithm 1 Double DQN with proportional prioritization

Input: minibatch k, step-size η, replay period K and size N, exponents α and β, budget T.
Initialize replay memory $\mathcal{H} = \emptyset$, $\Delta = 0$, $p_1 = 1$
Observe S_0 and choose $A_0 \sim \pi_\theta(S_0)$
for $t = 1$ **to** T **do**
  Observe S_t, R_t, γ_t
  Store transition $(S_{t-1}, A_{t-1}, R_t, \gamma_t, S_t)$ in $\mathcal{H}$ with maximal priority $p_t = \max_{i<t} p_i$
  if $t \equiv 0 \mod K$ **then**
    for $j = 1$ **to** k **do**
      Sample transition $j \sim P(j) = p_j^\alpha / \sum_i p_i^\alpha$
      Compute importance-sampling weight $w_j = (N \cdot P(j))^{-\beta} / \max_i w_i$
      Compute TD-error $\delta_j = R_j + \gamma_j Q_{\text{target}}(S_j, \arg\max_a Q(S_j, a)) - Q(S_{j-1}, A_{j-1})$
      Update transition priority $p_j \leftarrow |\delta_j|$
      Accumulate weight-change $\Delta \leftarrow \Delta + w_j \cdot \delta_j \cdot \nabla_\theta Q(S_{j-1}, A_{j-1})$
    end for
    Update weights $\theta \leftarrow \theta + \eta \cdot \Delta$, reset $\Delta = 0$
    From time to time copy weights into target network $\theta_{\text{target}} \leftarrow \theta$
  end if
  Choose action $A_t \sim \pi_\theta(S_t)$
end for

Algorithm 1 of "Prioritized Experience Replay" by Tom Schaul et al.

Dueling Networks

Instead of computing directly $Q(s, a; \boldsymbol{\theta})$, we compose it from the following quantities:

- average return in a given state s, $V(s; \boldsymbol{\theta}) = \frac{1}{|\mathcal{A}|} \sum_a Q(s, a; \boldsymbol{\theta})$,
- advantage function computing an **advantage** $Q(s, a; \boldsymbol{\theta}) - V(s; \theta)$ of action a in state s.

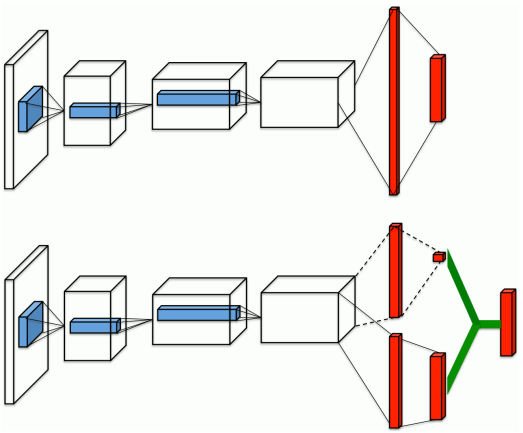


Figure 1 of "Dueling Network Architectures for Deep Reinforcement Learning" by Ziyu Wang et al.

$$Q(s, a) \stackrel{\text{def}}{=} V\big(f(s; \zeta); \eta\big) + A\big(f(s; \zeta), a; \psi\big) - \frac{\sum_{a' \in \mathcal{A}} A(f(s; \zeta), a'; \psi)}{|\mathcal{A}|}$$

Dueling Networks

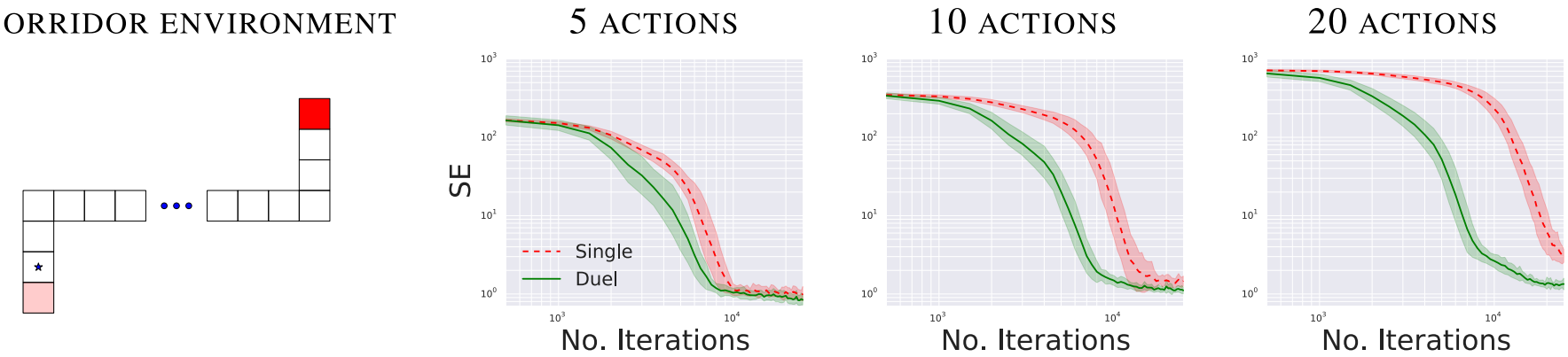


Figure 3. **(a)** The corridor environment. The star marks the starting state. The redness of a state signifies the reward the agent receives upon arrival. The game terminates upon reaching either reward state. The agent's actions are going up, down, left, right and no action. Plots **(b), (c)** and **(d)** shows squared error for policy evaluation with 5, 10, and 20 actions on a log-log scale. The dueling network (Duel) consistently outperforms a conventional single-stream network (Single), with the performance gap increasing with the number of actions.

Figure 3 of "Dueling Network Architectures for Deep Reinforcement Learning" by Ziyu Wang et al.

Dueling Networks

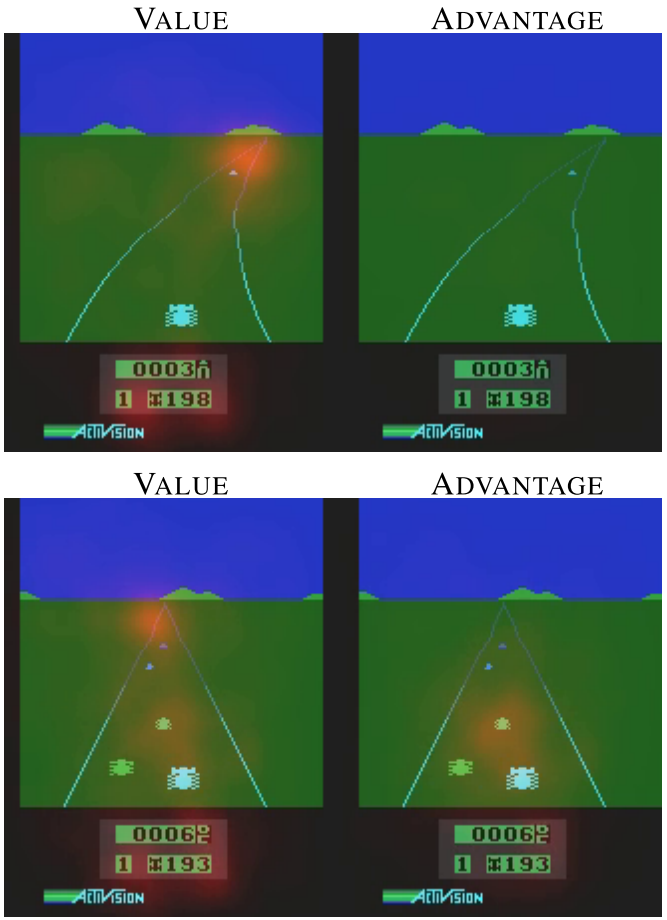


Figure 2 of "Dueling Network Architectures for Deep Reinforcement Learning" by Ziyu Wang et al.

Dueling Networks

Results on all 57 games (retraining the original DQN on the 8 missing games). `Single` refers to DDQN with a direct computation of $Q(s, a; \boldsymbol{\theta})$, `Clip` corresponds to additional gradient clipping to norm at most 10 and larger first hidden layer (so that duelling and single have roughly the same number of parameters).

	30 no-ops		Human Starts	
	Mean	**Median**	**Mean**	**Median**
Prior. Duel Clip	**591.9%**	**172.1%**	**567.0%**	**115.3%**
Prior. Single	434.6%	123.7%	386.7%	112.9%
Duel Clip	**373.1%**	**151.5%**	**343.8%**	**117.1%**
Single Clip	341.2%	132.6%	302.8%	114.1%
Single	307.3%	117.8%	332.9%	110.9%
Nature DQN	227.9%	79.1%	219.6%	68.5%

Table 1 of "Dueling Network Architectures for Deep Reinforcement Learning" by Ziyu Wang et al.