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Deep Reinforcement Learning

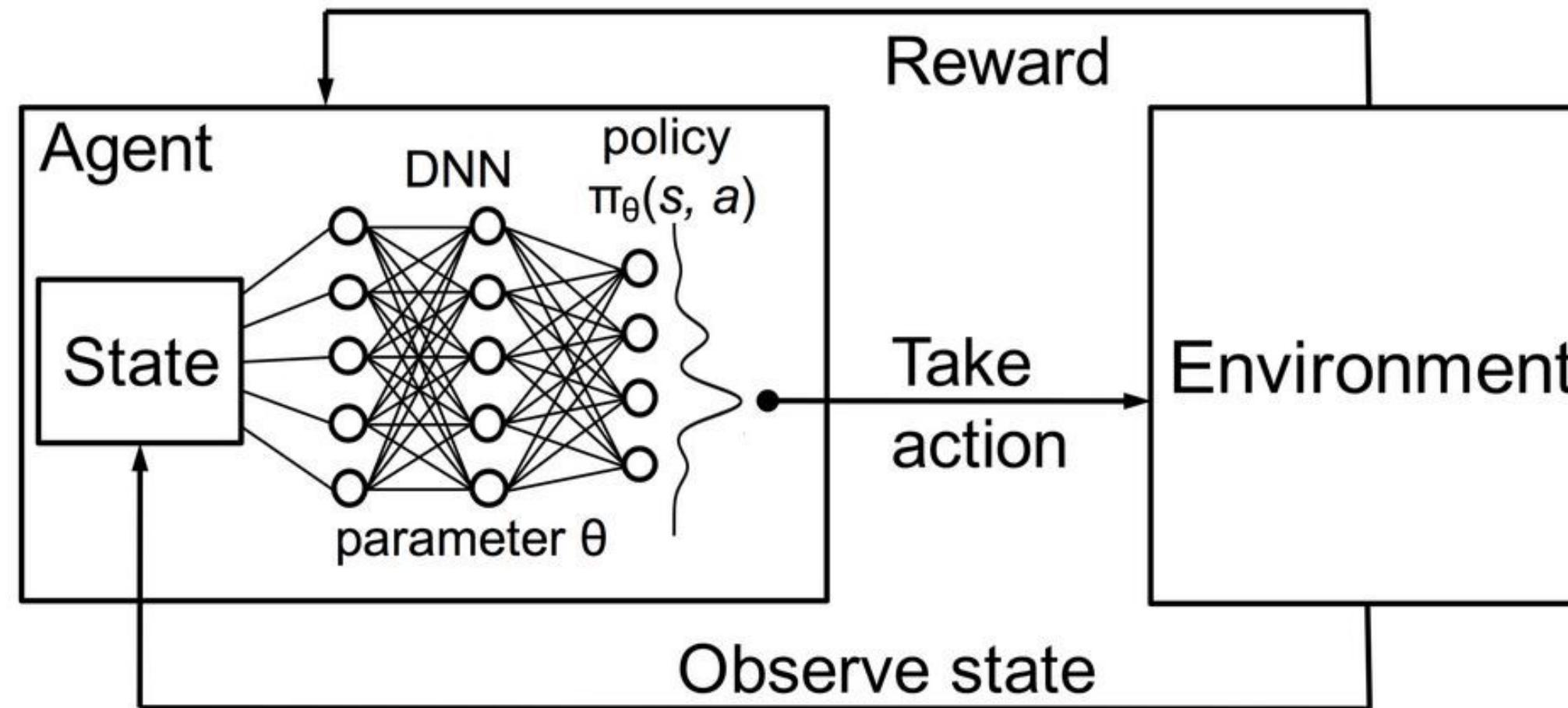
Deep Q-Learning

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<https://tu-chemnitz.de/informatik/KI/edu/deeprl>

Value-based deep RL



- The basic idea in **value-based deep RL** is to approximate the Q-values in each possible state, using a **deep neural network** with free parameters θ :

$$Q_\theta(s, a) \approx Q^\pi(s, a) = \mathbb{E}_\pi(R_t | s_t = s, a_t = a)$$

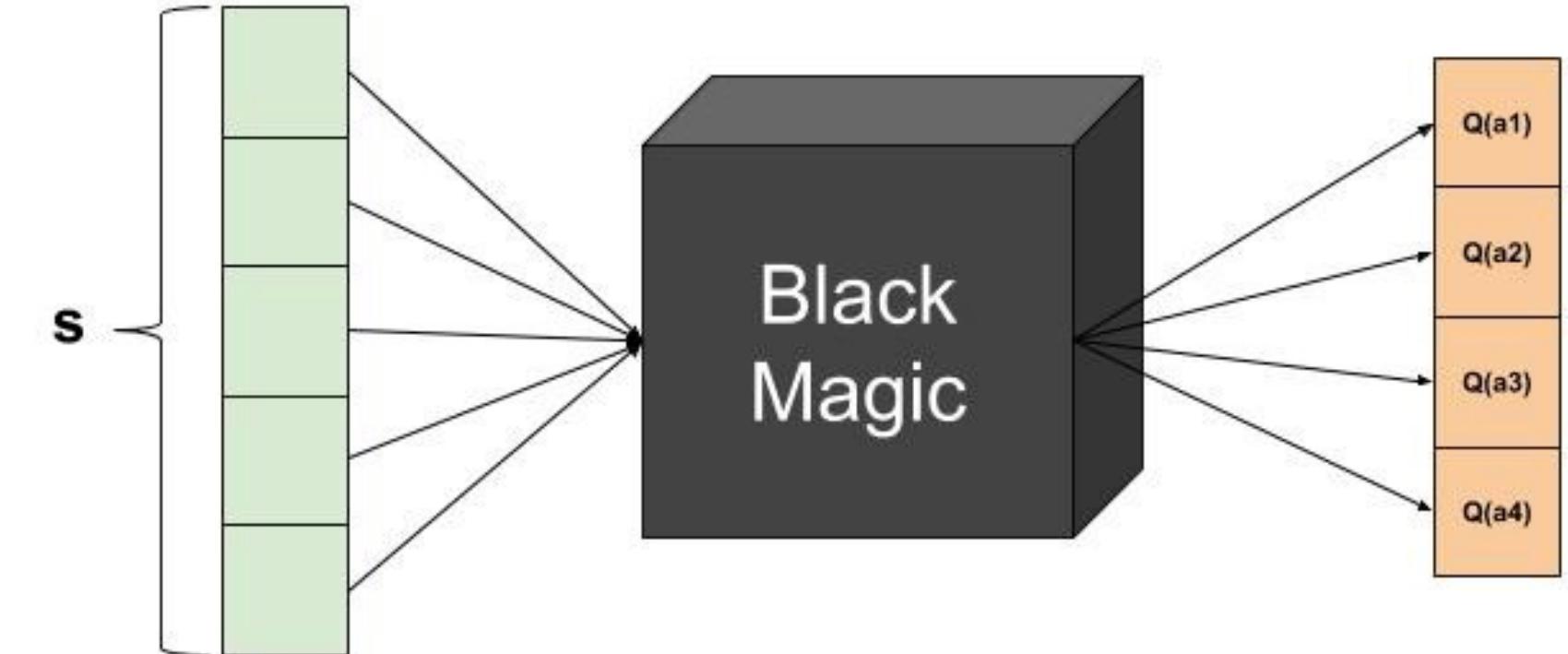
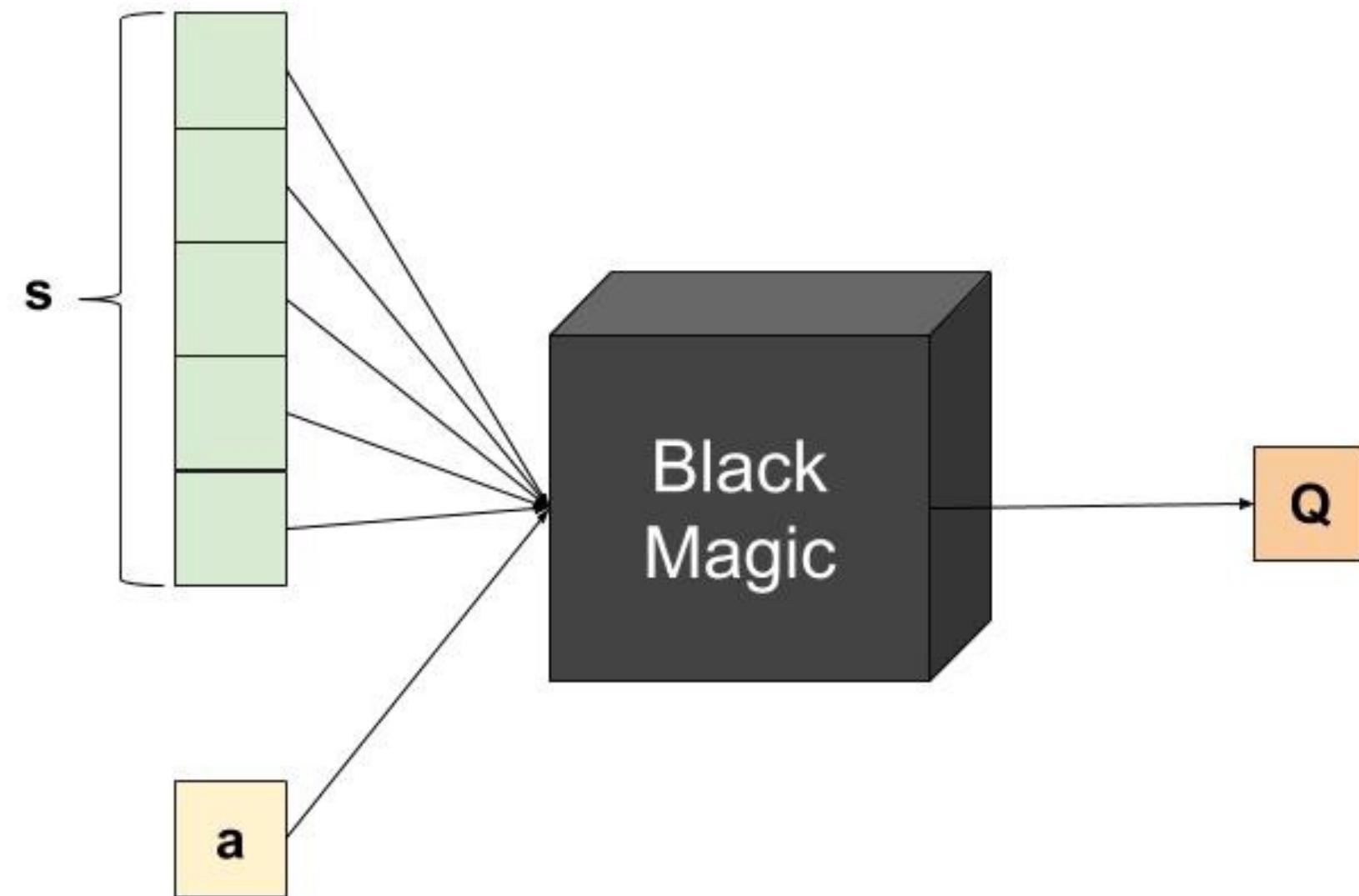
- The Q-values now depend on the parameters θ of the DNN.
- The derived policy π_θ uses for example an ϵ -greedy or softmax action selection scheme over the estimated Q-values:

$$\pi_\theta(s, a) \leftarrow \text{Softmax}(Q_\theta(s, a))$$

Function approximators to learn the Q-values

There are two possibilities to approximate Q-values $Q_\theta(s, a)$:

- The DNN approximates the Q-value of a single (s, a) pair.
- The DNN approximates the Q-value of all actions a in a state s .



- The action space must be discrete (one neuron per action).
- The action space can be continuous.

First naive approach: Q-learning with function approximation

We could simply adapt Q-learning with FA to the DNN:

- Initialize the deep neural network with parameters θ .
- Start from an initial state s_0 .
- for $t \in [0, T_{\text{total}}]$:
 - Select a_t using a softmax over the Q-values $Q_\theta(s_t, a)$.
 - Take a_t , observe r_{t+1} and s_{t+1} .
 - Update the parameters θ by minimizing the loss function:

$$\mathcal{L}(\theta) = (r_{t+1} + \gamma \max_{a'} Q_\theta(s_{t+1}, a') - Q_\theta(s_t, a_t))^2$$

- if s_t is terminal: sample another initial state s_0 .

Remark: We will now omit the break for terminal states, it is always implicitly here.

DNN need stochastic gradient descent

- This naive approach will not work: DNNs cannot learn from single examples (online learning = instability).
- DNNs require **stochastic gradient descent** (SGD):

$$\mathcal{L}(\theta) = E_{\mathcal{D}}(||\mathbf{t} - \mathbf{y}||^2) \approx \frac{1}{K} \sum_{i=1}^K ||\mathbf{t}_i - \mathbf{y}_i||^2$$

- The loss function is estimated by **sampling** a minibatch of K **i.i.d** samples from the training set to compute the loss function and update the parameters θ .
- This is necessary to avoid local minima of the loss function.
- Although Q-learning can learn from single transitions, it is not possible using DNN.
- Why not using the last K transitions to train the network? We could store them in a **transition buffer** and train the network on it.

Second naive approach: Q-learning with a transition buffer

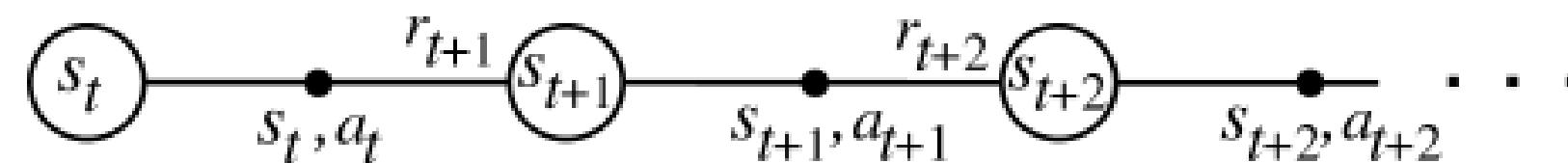
- Initialize the deep neural network with parameters θ .
- Initialize an empty **transition buffer** \mathcal{D} of size K : $\{(s_k, a_k, r_k, s'_k)\}_{k=1}^K$.
- for $t \in [0, T_{\text{total}}]$:
 - Select a_t using a softmax over the Q-values $Q_\theta(s_t, a)$.
 - Take a_t , observe r_{t+1} and s_{t+1} .
 - Store $(s_t, a_t, r_{t+1}, s_{t+1})$ in the transition buffer.
 - Every K steps:
 - Update the parameters θ using the transition buffer:

$$\mathcal{L}(\theta) = \frac{1}{K} \sum_{k=1}^K (r_k + \gamma \max_{a'} Q_\theta(s'_k, a') - Q_\theta(s_k, a_k))^2$$

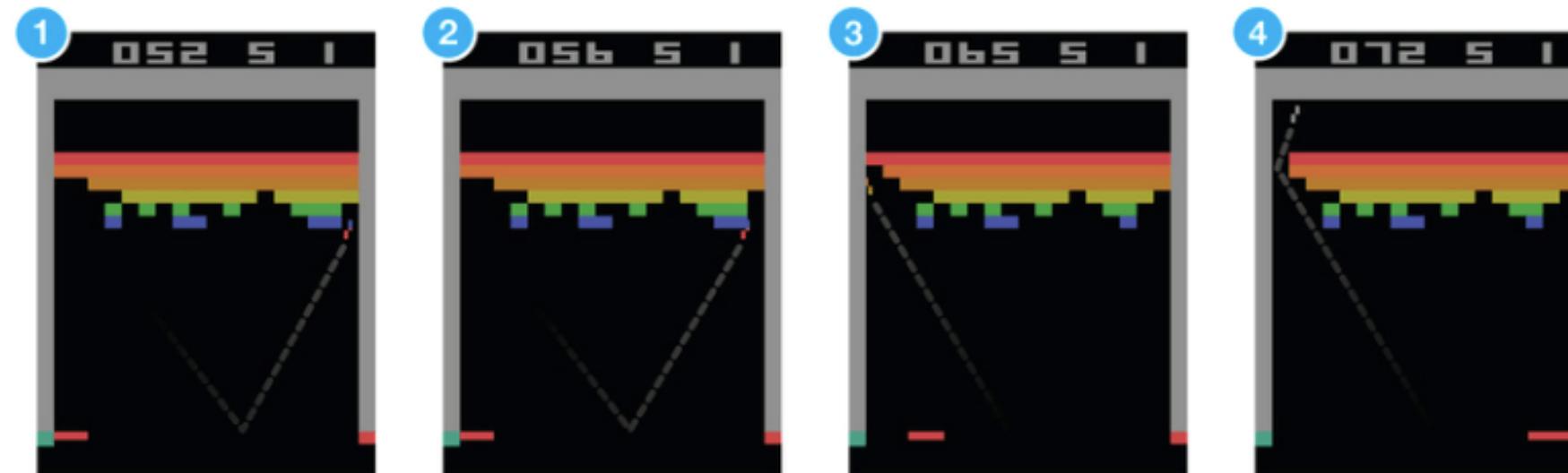
- Empty the transition buffer.

Correlated inputs

- Unfortunately, this does not work either.
- The last K transitions (s, a, r, s') are not **i.i.d** (independent and identically distributed).



- The transition $(s_{t+1}, a_{t+1}, r_{t+2}, s_{t+2})$ **depends** on $(s_t, a_t, r_{t+1}, s_{t+1})$ by definition, i.e. the transitions are **correlated**.
- Even worse, when playing video games, successive frames will be very similar or even identical.



- The actions are also correlated: you move the paddle to the left for several successive steps.

Correlated inputs

- Feeding transitions sequentially to a DNN is the same as giving all MNIST 0's to a DNN, then all 1's, etc... It does not work.



Sequential-Correlated



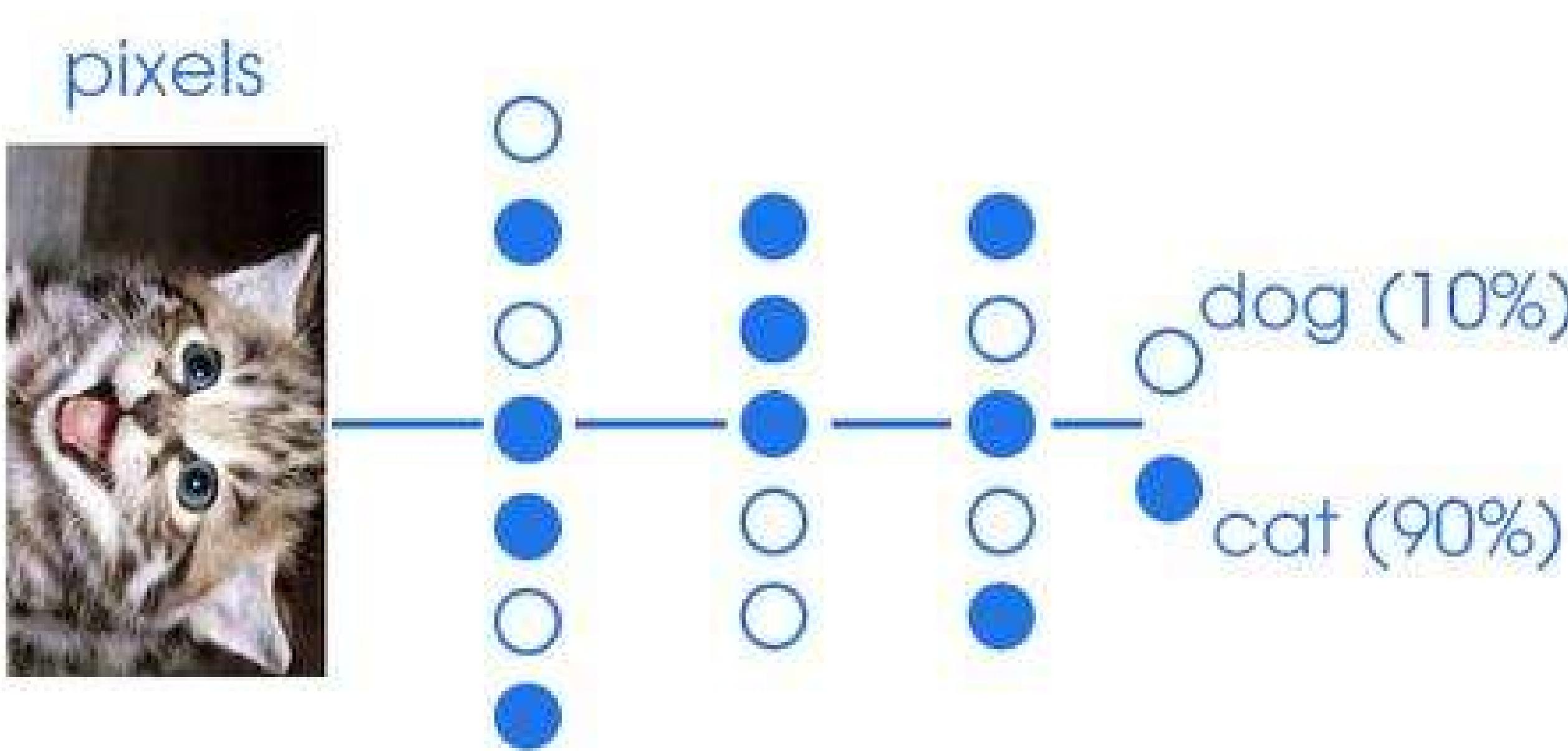
Varied Data Dist.

- In SL, we have all the training data **before** training: it is possible to get i.i.d samples by shuffling the training set between two epochs.
- In RL, we create the “training set” (transitions) **during** training: the samples are not i.i.d as we act sequentially over time.

Non-stationarity

- In SL, the **targets t** do not change over time: an image of a cat stays an image of a cat throughout learning.

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathbf{x}, \mathbf{t} \sim \mathcal{D}} [||\mathbf{t} - F_\theta(\mathbf{x})||^2]$$



- The problem is said **stationary**, as the distribution of the data does not change over time.

Non-stationarity

- In RL, the **targets** $t = r + \gamma \max_{a'} Q_\theta(s', a')$ do change over time:
 - $Q_\theta(s', a')$ depends on θ , so after one optimization step, all targets have changed!
 - As we improve the policy over training, we collect higher returns.

$$\mathcal{L}(\theta) = \mathbb{E}_{s,a \sim \pi_\theta} [(r + \gamma \max_{a'} Q_\theta(s', a') - Q_\theta(s, a))^2]$$

- NN do not like this. After a while, they give up and settle on a **suboptimal** policy.

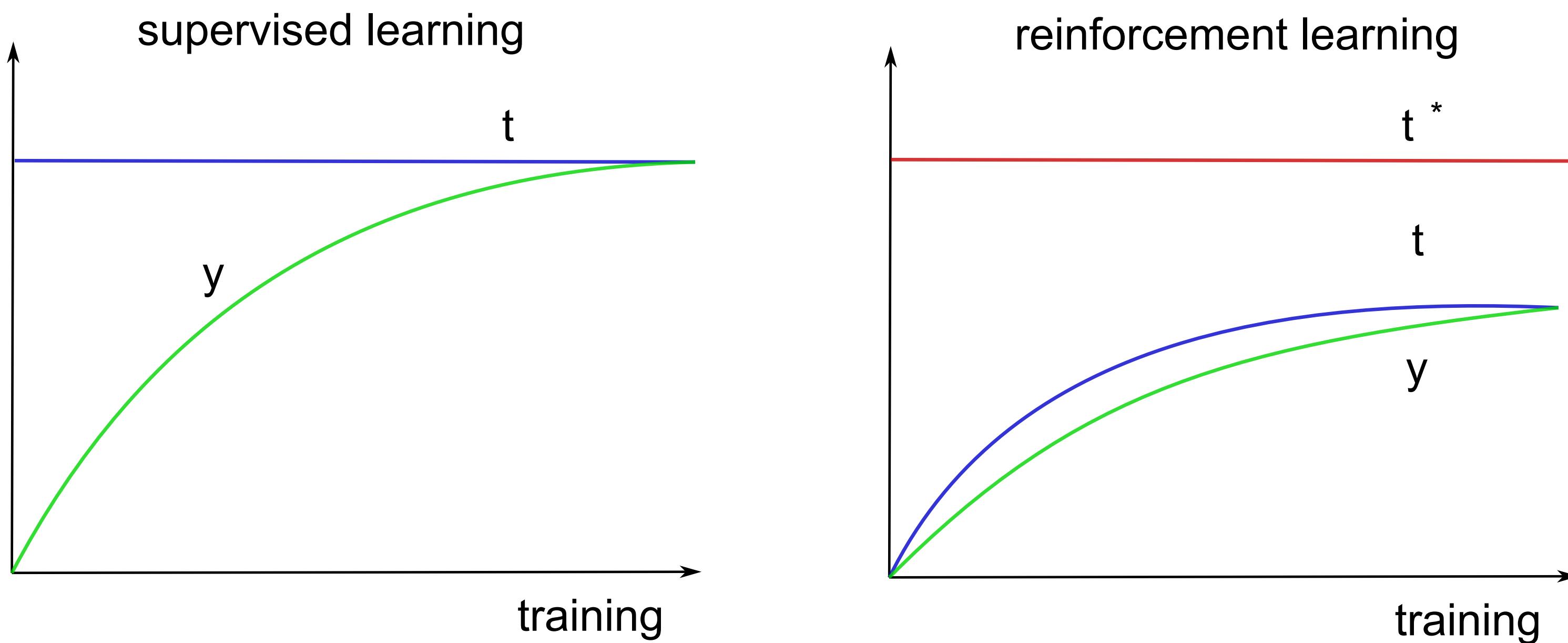
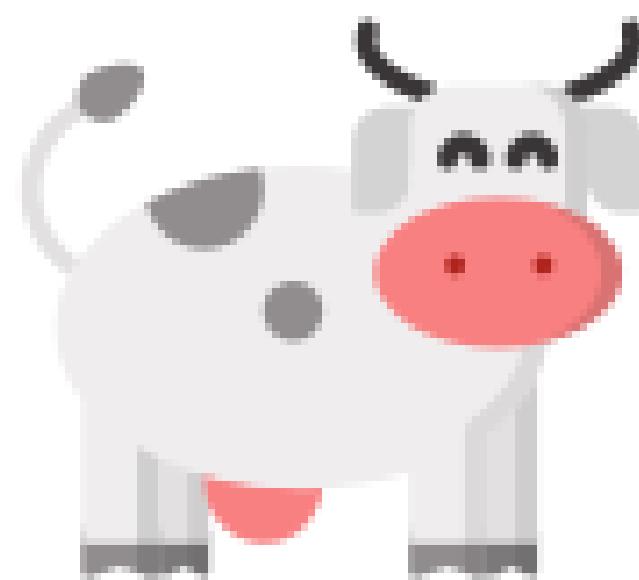
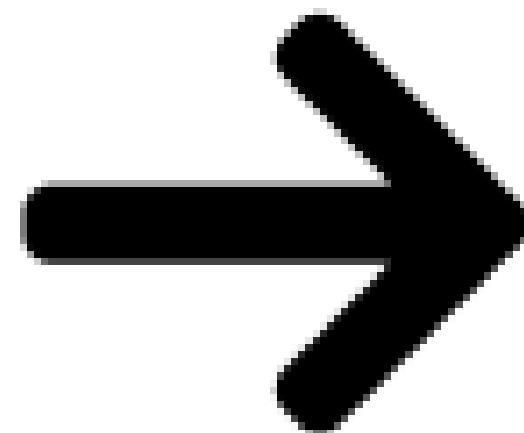


Illustration of non-stationary targets

- We want our value estimates to “catch” the true values.



Q estimation



Q target

Illustration of non-stationary targets

- We update our estimate to come closer to the target.

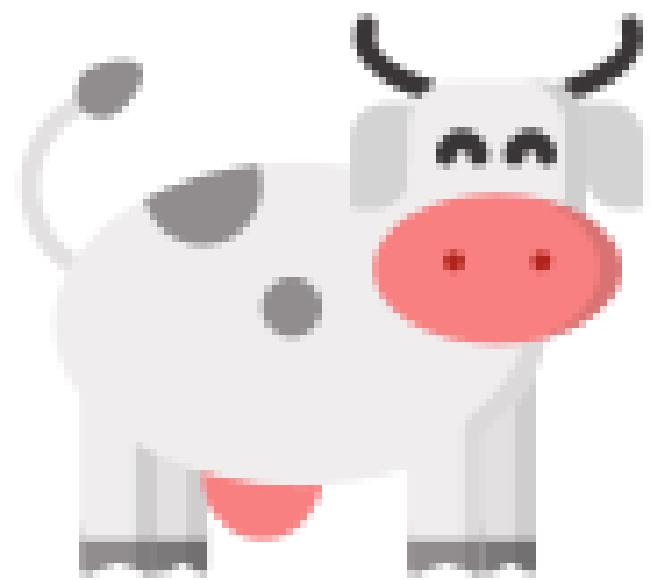
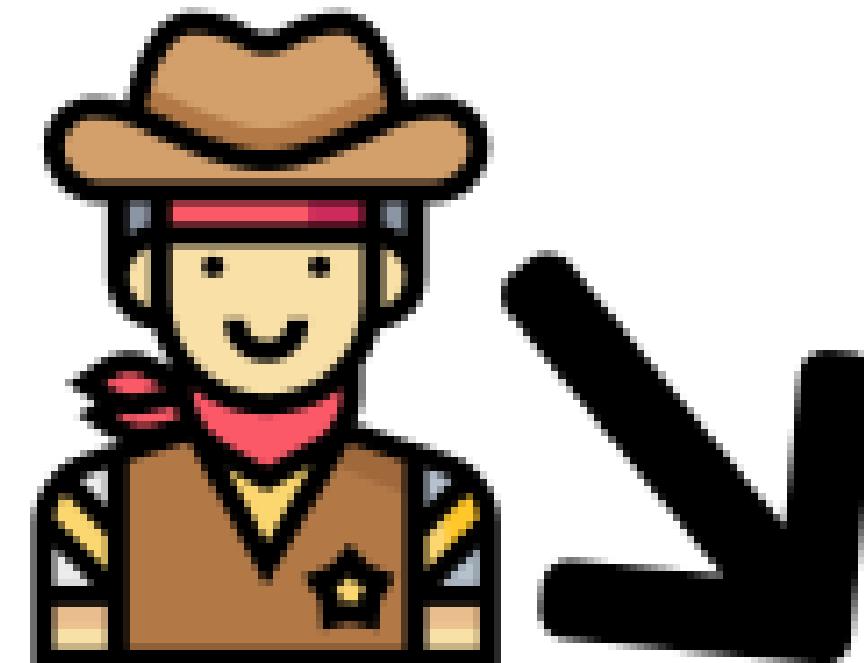


Illustration of non-stationary targets

- But the target moves! We need to update again.

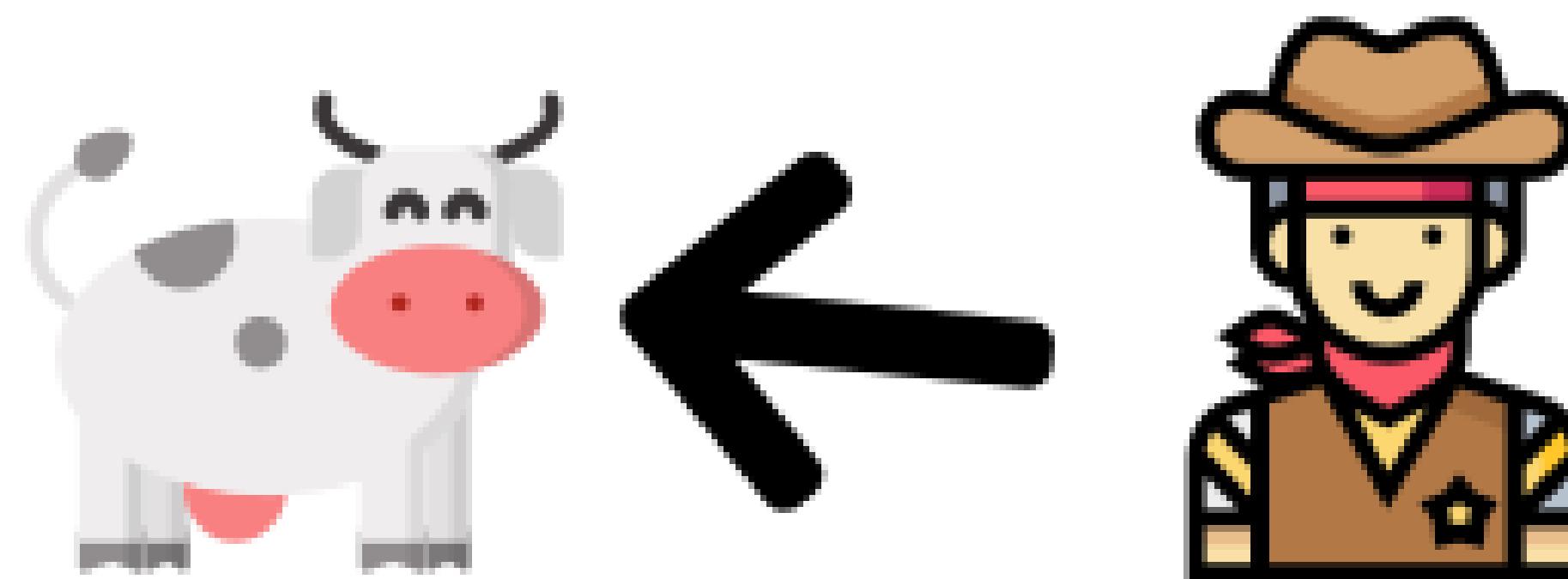
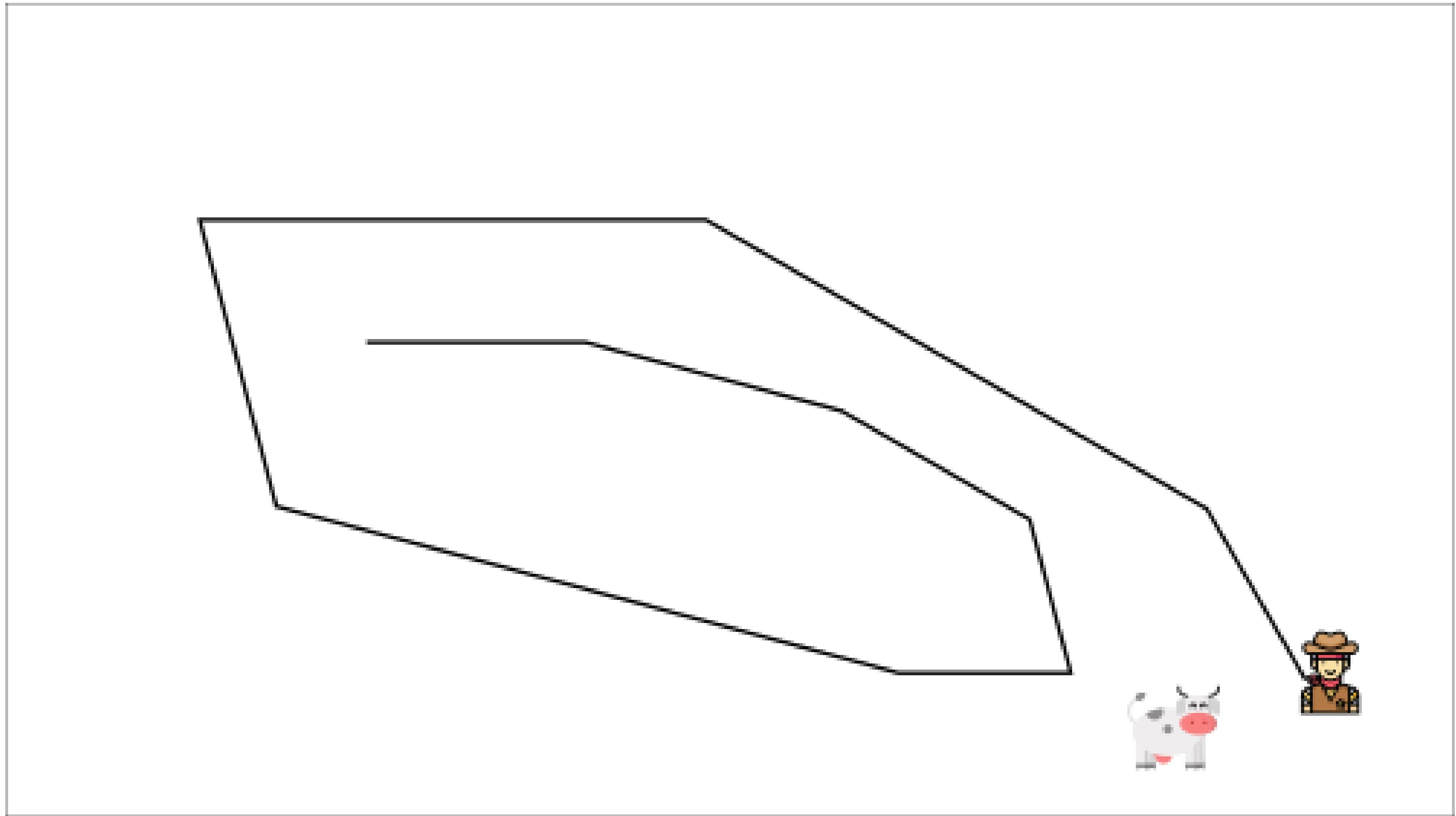


Illustration of non-stationary targets

- This leads to very strange and inefficient optimization paths



1 - Deep Q-networks (DQN)

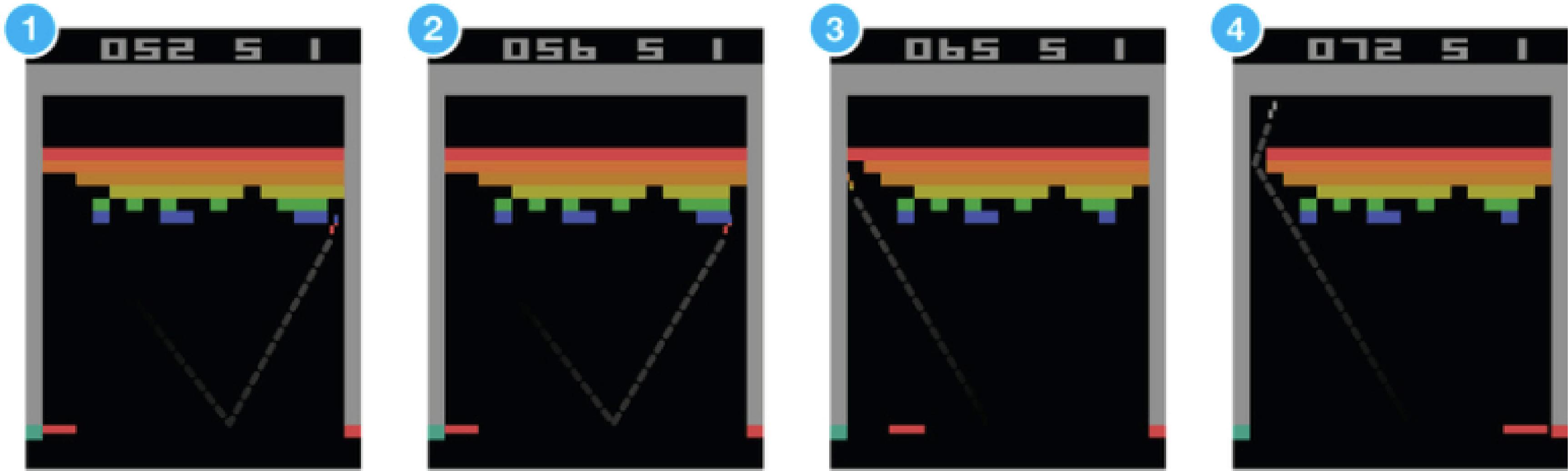
Playing Atari with Deep Reinforcement Learning

Volodymyr Mnih Koray Kavukcuoglu David Silver Alex Graves Ioannis Antonoglou

Daan Wierstra Martin Riedmiller

DeepMind Technologies

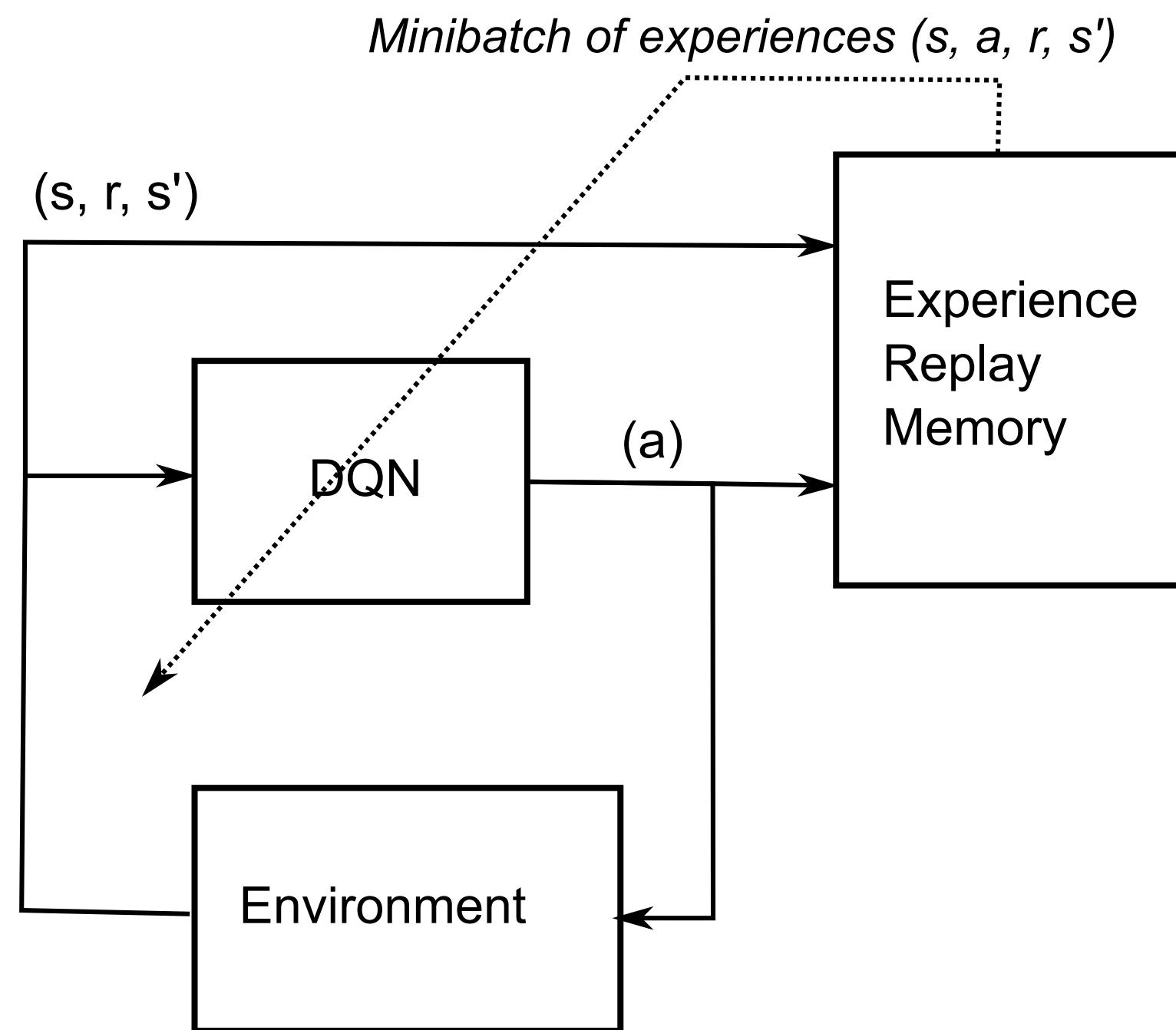
Problem with non-linear approximators and RL



- Non-linear approximators never really worked with RL before 2013 because of:
 1. The correlation between successive inputs or outputs.
 2. The non-stationarity of the problem.
- These two problems are very bad for deep networks, which end up overfitting the learned episodes or not learning anything at all.
- Deepmind researchers proposed to use two classical ML tricks to overcome these problems:
 1. experience replay memory.
 2. target networks.

Experience replay memory

- To avoid correlation between samples, (Mnih et al. 2015) proposed to store the (s, a, r, s') transitions in a huge **experience replay memory** or **replay buffer** \mathcal{D} (e.g. 1 million transitions).



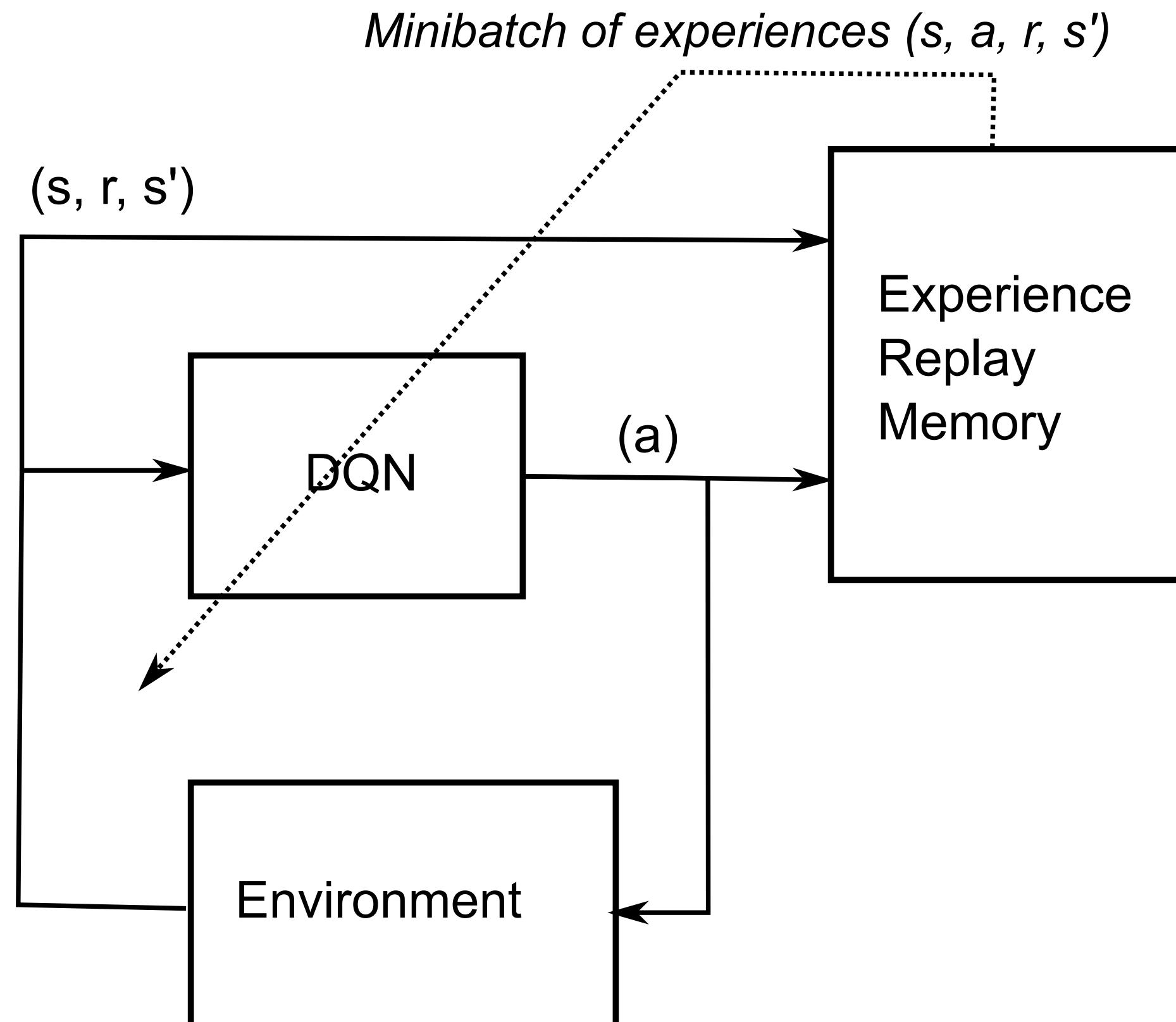
- When the buffer is full, we simply overwrite old transitions.
- The Q-learning update is only applied on a **random minibatch** of those past experiences, not the last transitions.
- This ensures the independence of the samples (non-correlated samples).

Experience replay memory

- Initialize value network Q_θ .
- Initialize experience replay memory \mathcal{D} of maximal size N .
- for $t \in [0, T_{\text{total}}]$:
 - Select an action a_t based on $Q_\theta(s_t, a)$, observe s_{t+1} and r_{t+1} .
 - Store $(s_t, a_t, r_{t+1}, s_{t+1})$ in the experience replay memory.
 - Every T_{train} steps:
 - Sample a minibatch \mathcal{D}_s randomly from \mathcal{D} .
 - For each transition (s_k, a_k, r_k, s'_k) in the minibatch:
 - Compute the target value $t_k = r_k + \gamma \max_{a'} Q_\theta(s'_k, a')$
 - Update the value network Q_θ on \mathcal{D}_s to minimize:

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathcal{D}_s} [(t_k - Q_\theta(s_k, a_k))^2]$$

Experience replay memory



- But wait! The samples of the minibatch are still not i.i.d, as they are not **identically distributed**:
 - Some samples were generated with a very old policy π_{θ_0} .
 - Some samples have been generated recently by the current policy π_{θ} .
- The samples of the minibatch do not come from the same distribution, so this should not work.

Experience replay memory

- This should not work, except if you use an **off-policy** algorithm, such as Q-learning!

$$Q^\pi(s, a) = \mathbb{E}_{s_t \sim \rho_b, a_t \sim b} [r_{t+1} + \gamma \max_a Q^\pi(s_{t+1}, a) | s_t = s, a_t = a]$$

- In Q-learning, you can take samples from **any** behavior policy b , as long as the coverage assumption stands:

$$\pi(s, a) > 0 \Rightarrow b(s, a) > 0$$

- Here, the behavior policy b is a kind of “superset” of all past policies π used to fill the ERM, so it “covers” the current policy.

$$b = \{\pi_{\theta_0}, \pi_{\theta_1}, \dots, \pi_{\theta_t}\}$$

- Samples from b are i.i.d, so Q-learning is going to work.
- Note: **it is not possible to use an experience replay memory with on-policy algorithms.**

$$Q^\pi(s, a) = \mathbb{E}_{s_t \sim \rho_\pi, a_t \sim \pi} [r_{t+1} + \gamma Q^\pi(s_{t+1}, a_{t+1}) | s_t = s, a_t = a]$$

- $a_{t+1} \sim \pi_\theta$ would not be the same between π_{θ_0} (which generated the sample) and π_{θ_t} (the current policy).

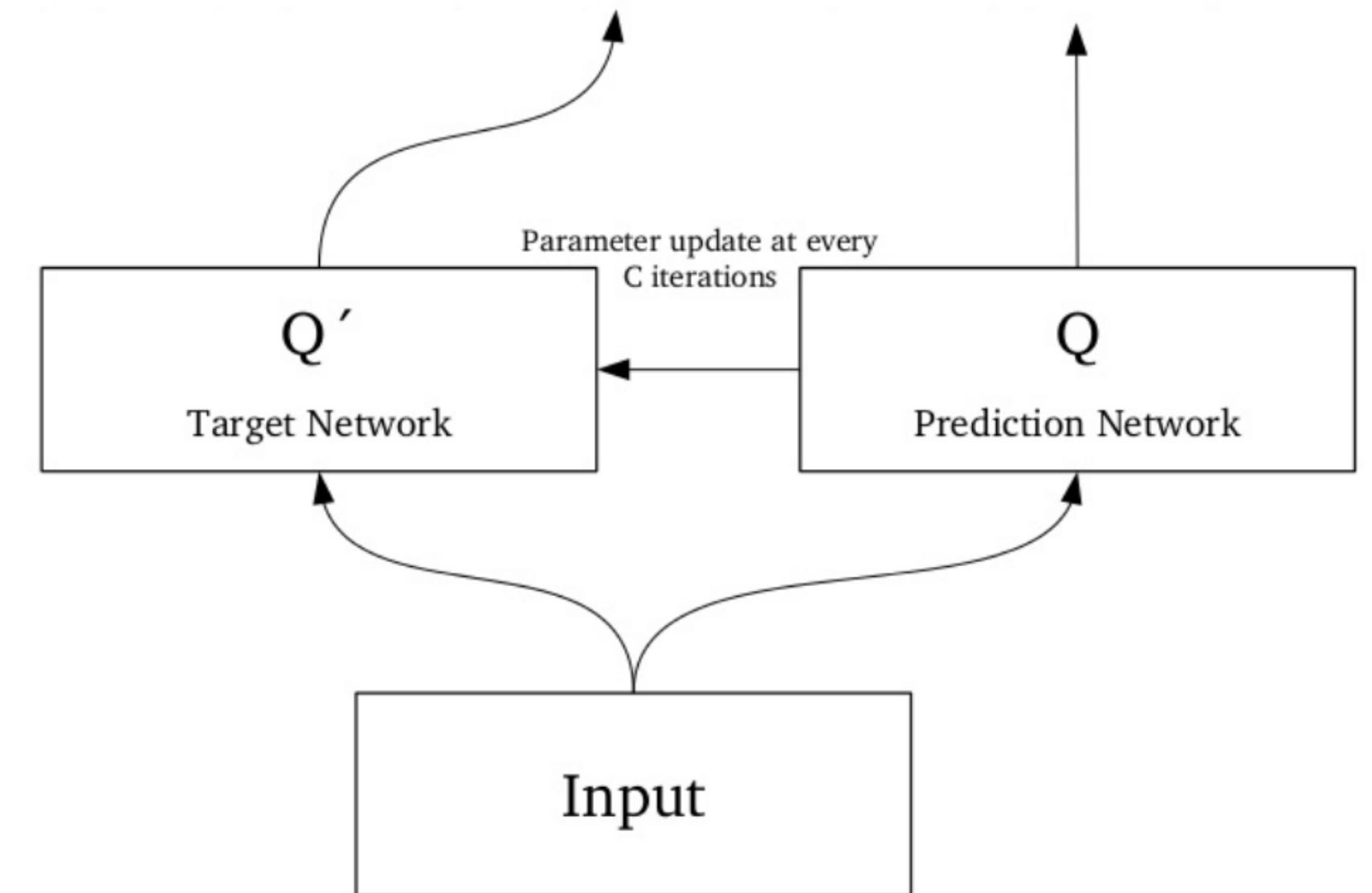
Target network

- The second problem when using DNN for RL is that the target is **non-stationary**, i.e. it changes over time: as the network becomes better, the Q-values have to increase.
- In DQN, the target for the update is not computed from the current deep network θ :

$$r + \gamma \max_{a'} Q_\theta(s', a')$$

but from a **target network** θ' updated only every few thousands of iterations.

$$r + \gamma \max_{a'} Q_{\theta'}(s', a')$$

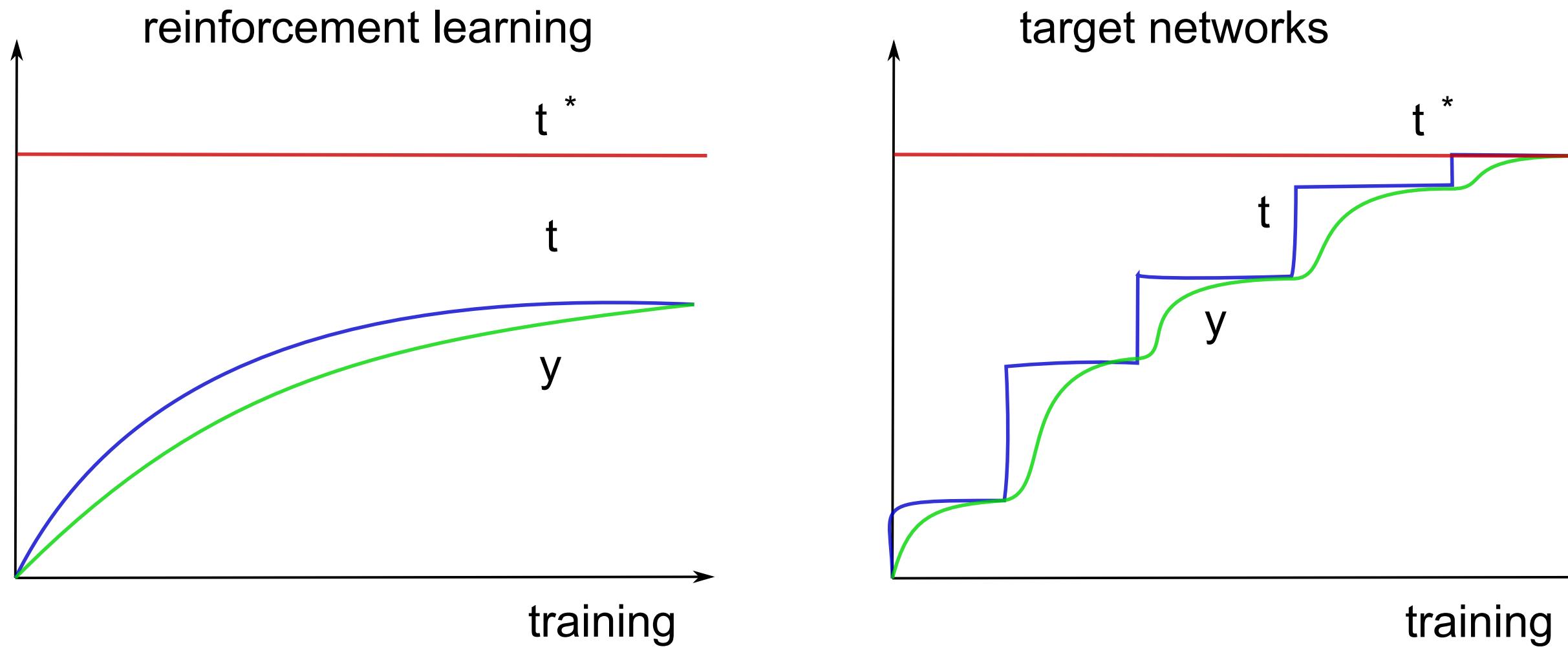


- θ' is simply a copy of θ from the past.
- DQN loss function:

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathcal{D}}[(r + \gamma \max_{a'} Q_{\theta'}(s', a')) - Q_\theta(s, a))^2]$$

Target network

- This allows the target $r + \gamma \max_{a'} Q_{\theta'}(s', a')$ to be **stationary** between two updates.
- It leaves time for the trained network to catch up with the targets.



- The update is simply replacing the parameters θ' with the trained parameters θ :

$$\theta' \leftarrow \theta$$

- The value network θ basically learns using an older version of itself...

DQN: Deep Q-network

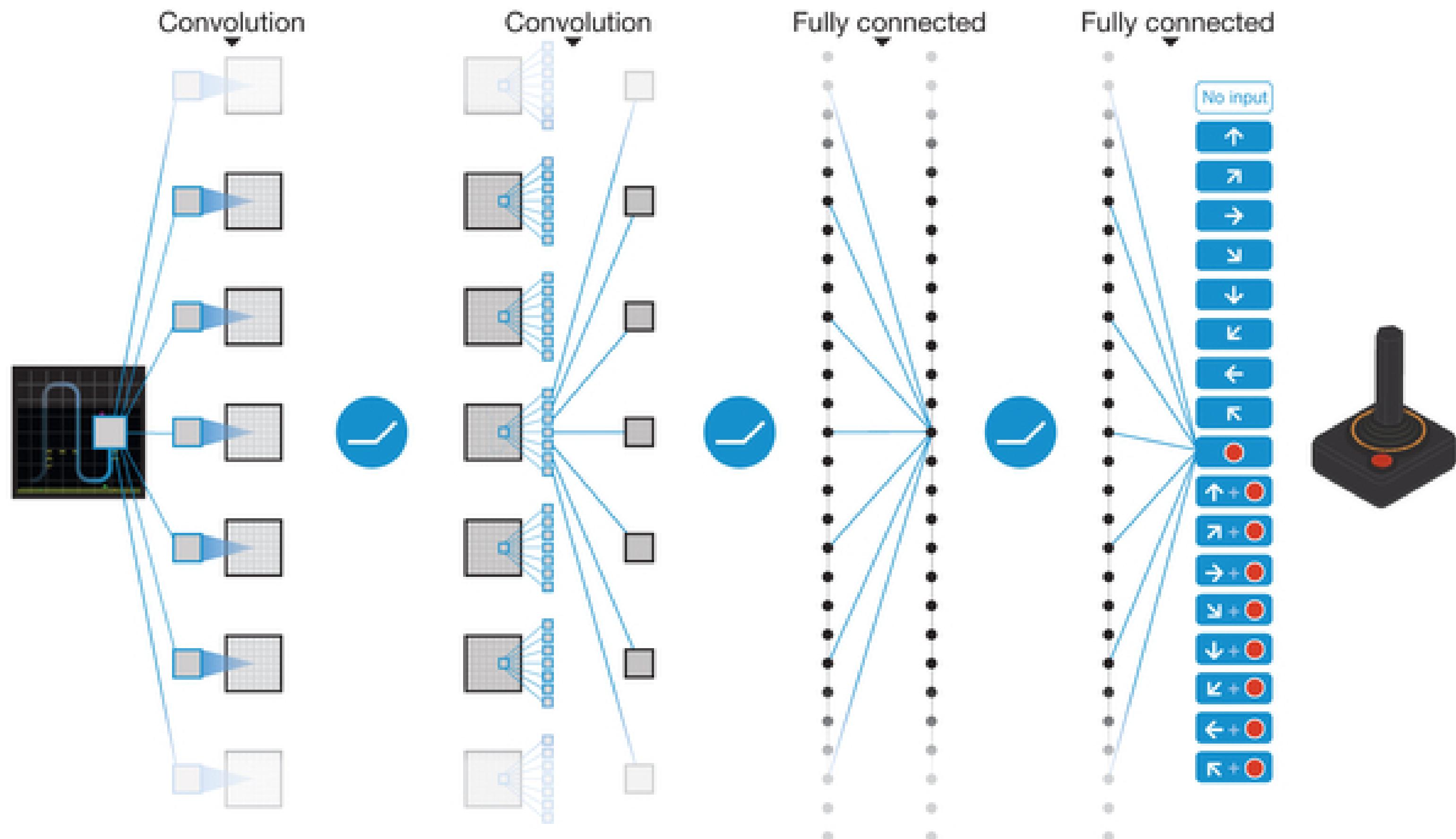
- Initialize value network Q_θ and target network $Q_{\theta'}$.
- Initialize experience replay memory \mathcal{D} of maximal size N .
- for $t \in [0, T_{\text{total}}]$:
 - Select an action a_t based on $Q_\theta(s_t, a)$, observe s_{t+1} and r_{t+1} .
 - Store $(s_t, a_t, r_{t+1}, s_{t+1})$ in the experience replay memory.
 - Every T_{train} steps:
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 - Update the value network Q_θ on \mathcal{D}_s to minimize:

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathcal{D}_s} [(t_k - Q_\theta(s_k, a_k))^2]$$

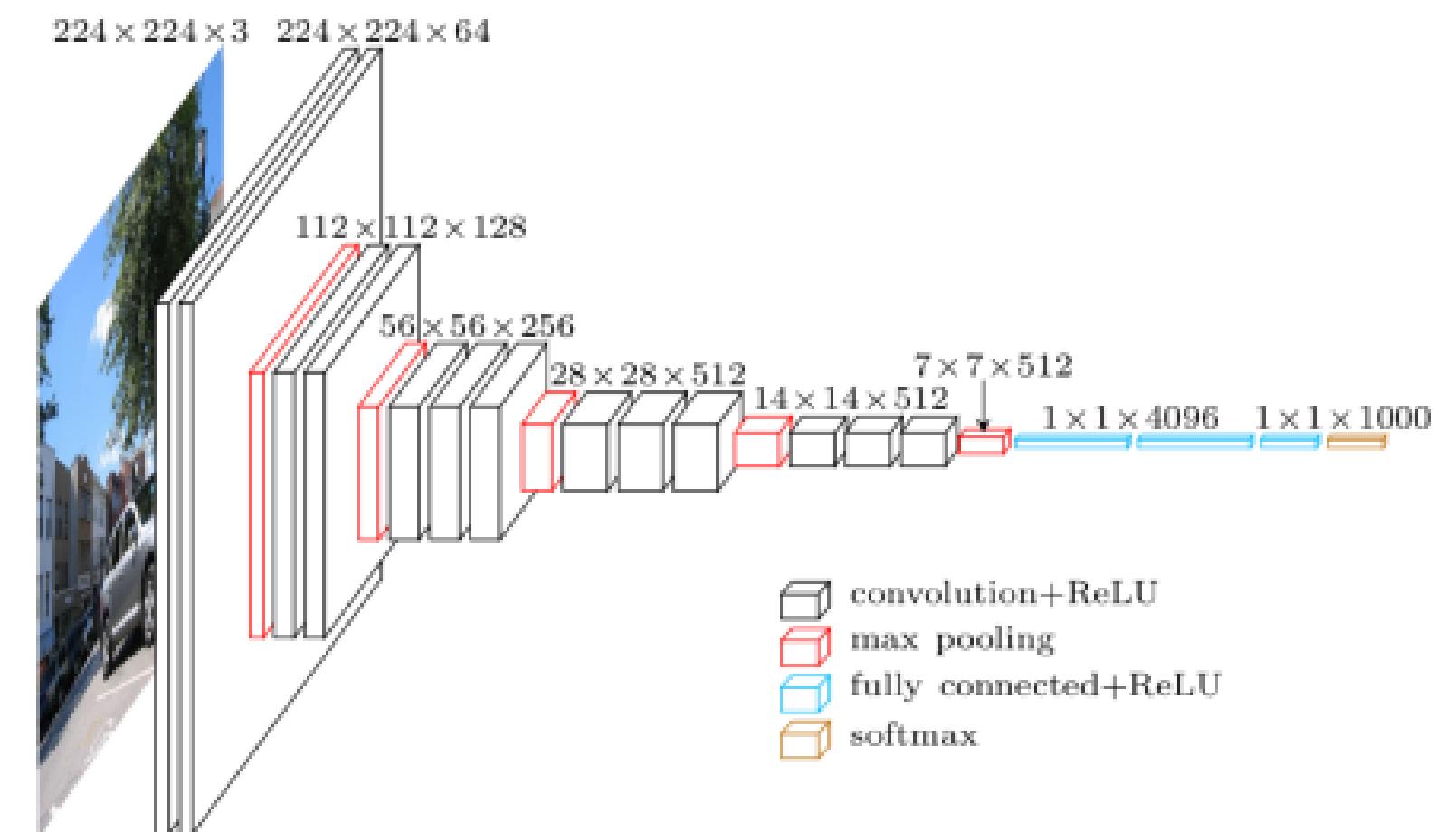
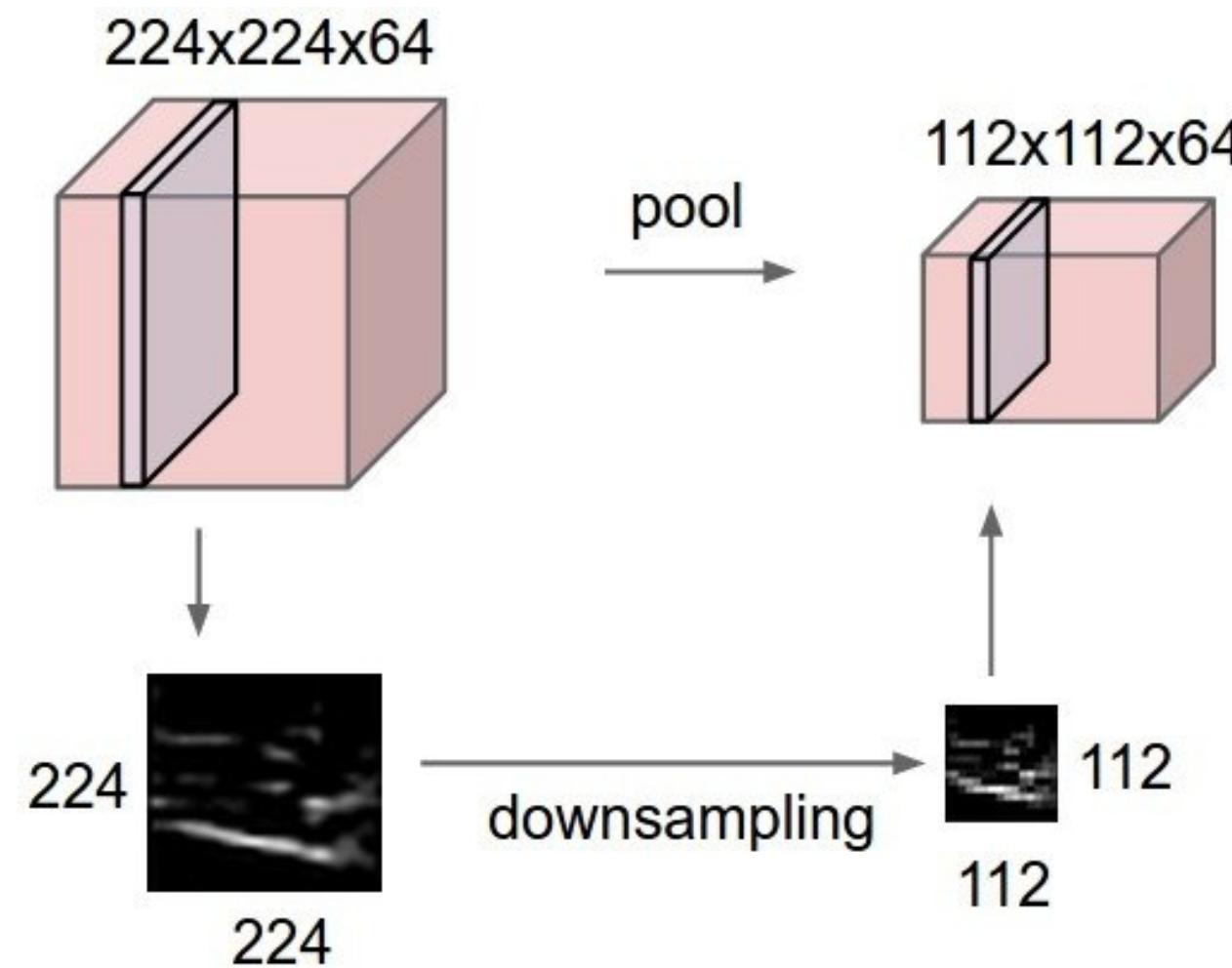
- Every T_{target} steps:
 - Update target network: $\theta' \leftarrow \theta$.

DQN: Deep Q-network

- The deep network can be anything. Deep RL is only about defining the loss function adequately.
- For pixel-based problems (e.g. video games), convolutional neural networks (without max-pooling) are the weapon of choice.



Why no max-pooling?



- The goal of max-pooling is to get rid of the spatial information in the image.
- For object recognition, you do not care whether the object is in the center or on the side of the image.
- Max-pooling brings **spatial invariance**.
- In video games, you **want** to keep the spatial information: the optimal action depends on where the ball is relative to the paddle.

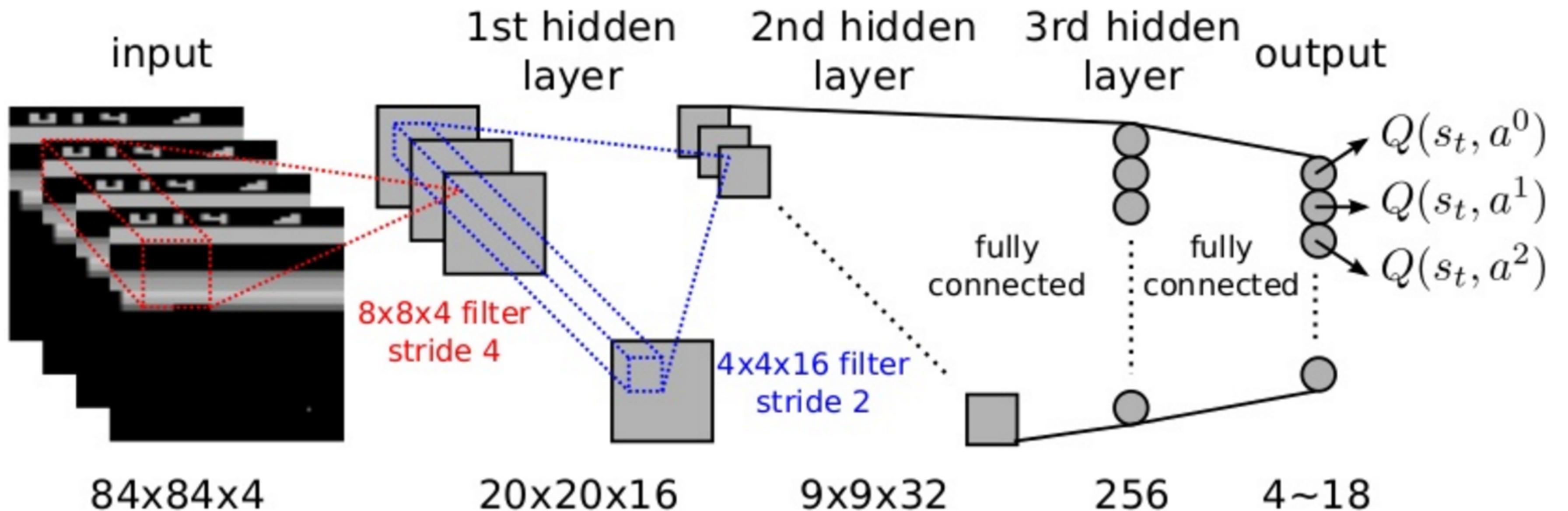
Are individual frames good representations of states?

- Is the ball moving from the child to the baseball player, or the other way around?



- Using video frames as states breaks the Markov property: the speed and direction of the ball is a very relevant information for the task, but not contained in a single frame.
- This characterizes a **Partially-observable Markov Decision Process** (POMDP).

Markov property in video games



- The simple solution retained in the original DQN paper is to **stack** the last four frames to form the state representation.
- Having the previous positions of the ball, the network can **learn** to infer its direction of movement.

DQN code in Keras

- Creating the CNN in keras / tensorflow / pytorch is straightforward:

```
1 model = Sequential()
2
3 model.add(Input((4, 84, 84)))
4
5 model.add(Conv2D(16, (8, 8), strides=(4, 4), activation='relu'))
6
7 model.add(Conv2D(32, (4, 4), strides=(2, 2), activation='relu'))
8
9 model.add(Flatten())
10
11 model.add(Dense(256, activation='relu'))
12
13 model.add(Dense(nb_actions, activation='linear'))
14
15 optimizer = RMSprop(lr=0.00025, rho=0.95, epsilon=0.01)
16
17 model.compile(optimizer, loss='mse')
```

DQN code in Keras

- Each step of the algorithm follows the GPI approach:

```
1 def q_iteration(env, model, state, memory):
2     # Choose the action with epsilon-greedy
3     if np.random.random() < epsilon:
4         action = env.action_space.sample()
5     else:
6         # Predict the Q-values for the current state and take the greedy action
7         values = model.predict([state])[0]
8         action = values.argmax()
9
10    # Play one game iteration
11    new_state, reward, done, _ = env.step(action)
12
13    # Append the transition to the replay buffer
14    memory.add(state, action, new_state, reward, done)
15
16    # Sample a minibatch from the memory and fit the DQN
17    s, a, r, s_, d = memory.sample_batch(32)
18    fit_batch(model, s, a, r, s_, d)
```

DQN code in Keras

- The only slight difficulty is actually to compute the targets for learning:

```
1 def fit_batch(model, states, actions, rewards, next_states, dones)
2
3     # Predict the Q-values in the current state
4     Q_values = model.predict(states)
5
6     # Predict the Q-values in the next state using the target model
7     next_Q_value = target_model.predict(next_states).max(axis=1)
8
9     # Terminal states have a value of 0
10    next_Q_value[dones] = 0.0
11
12    # Compute the target
13    targets = Q_values.copy()
14    for i in range(batch_size):
15        targets[i, actions[i]] = rewards[i] + self.gamma * next_Q_value[i]
16
17    # Train the model on the minibatch
18    self.model.fit(states, targets, epochs=1, batch_size=batch_size, verbose=0)
```

DQN training

- 50M frames (38 days of game experience) per game. Replay buffer of 1M frames.
- Action selection: ϵ -greedy with $\epsilon = 0.1$ and annealing. Optimizer: RMSprop with a batch size of 32.

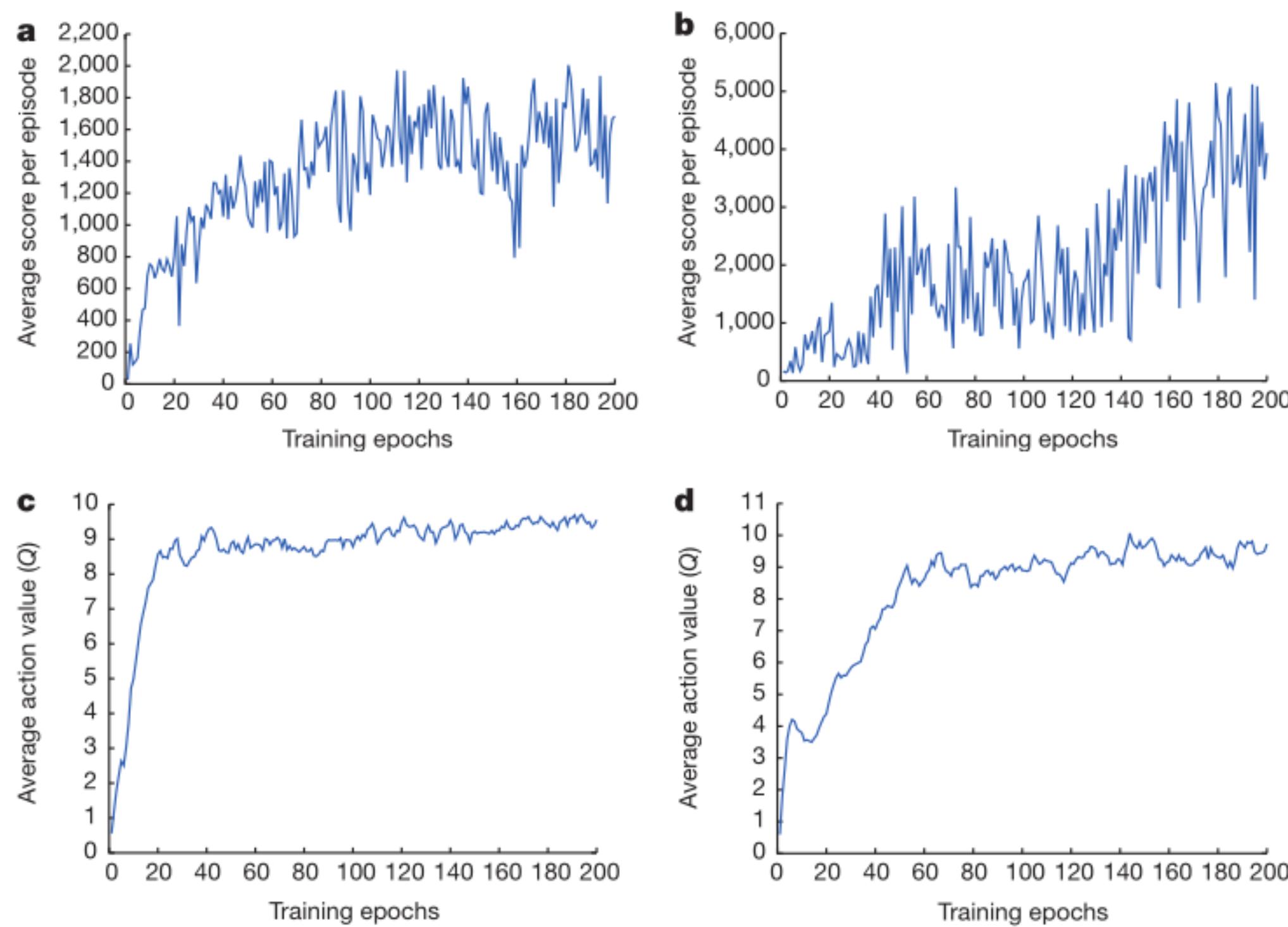


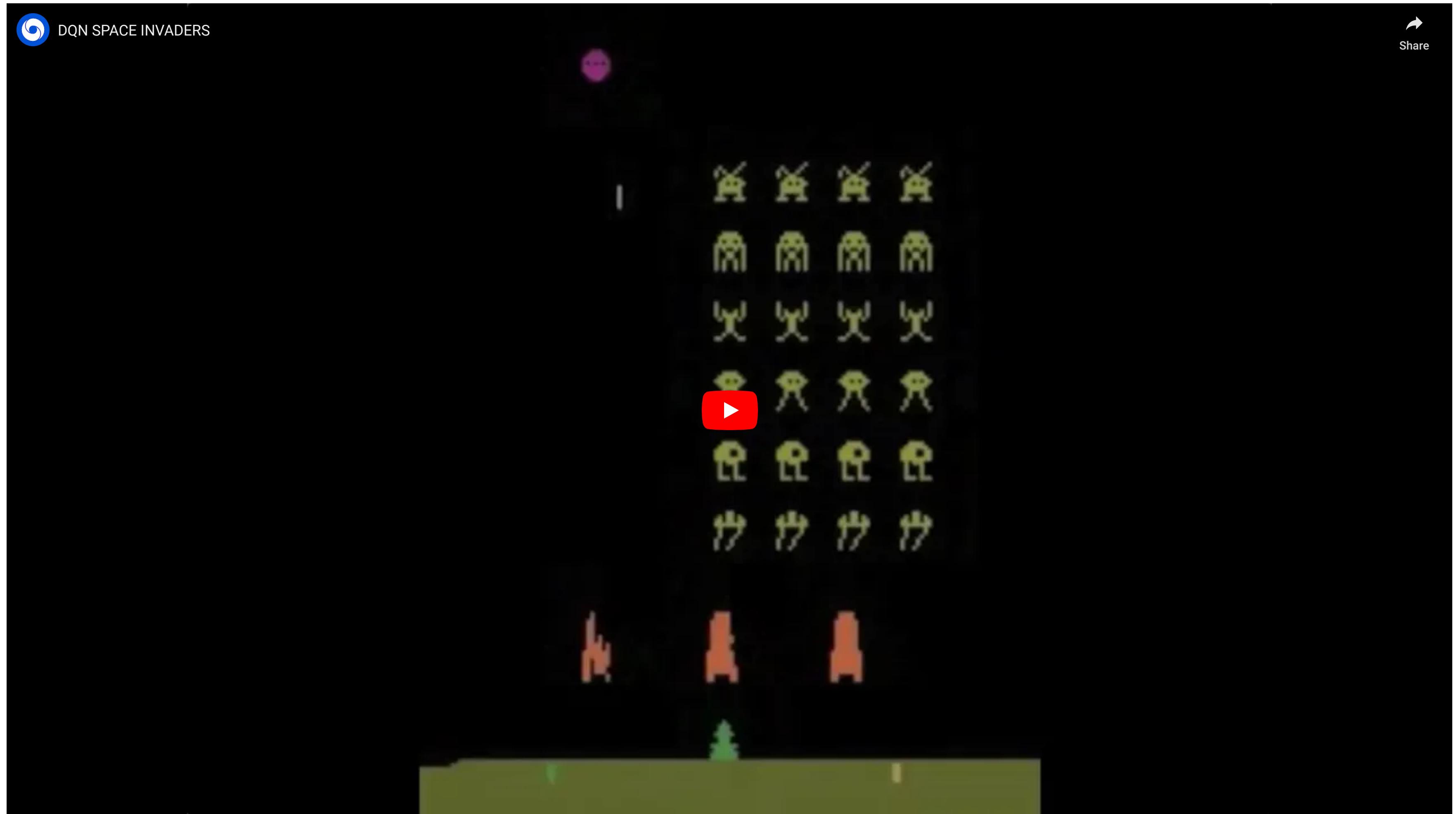
Figure 2 | Training curves tracking the agent’s average score and average predicted action-value. **a**, Each point is the average score achieved per episode after the agent is run with ϵ -greedy policy ($\epsilon = 0.05$) for 520 k frames on Space Invaders. **b**, Average score achieved per episode for Seaquest. **c**, Average predicted action-value on a held-out set of states on Space Invaders. Each point

on the curve is the average of the action-value Q computed over the held-out set of states. Note that Q-values are scaled due to clipping of rewards (see Methods). **d**, Average predicted action-value on Seaquest. See Supplementary Discussion for details.

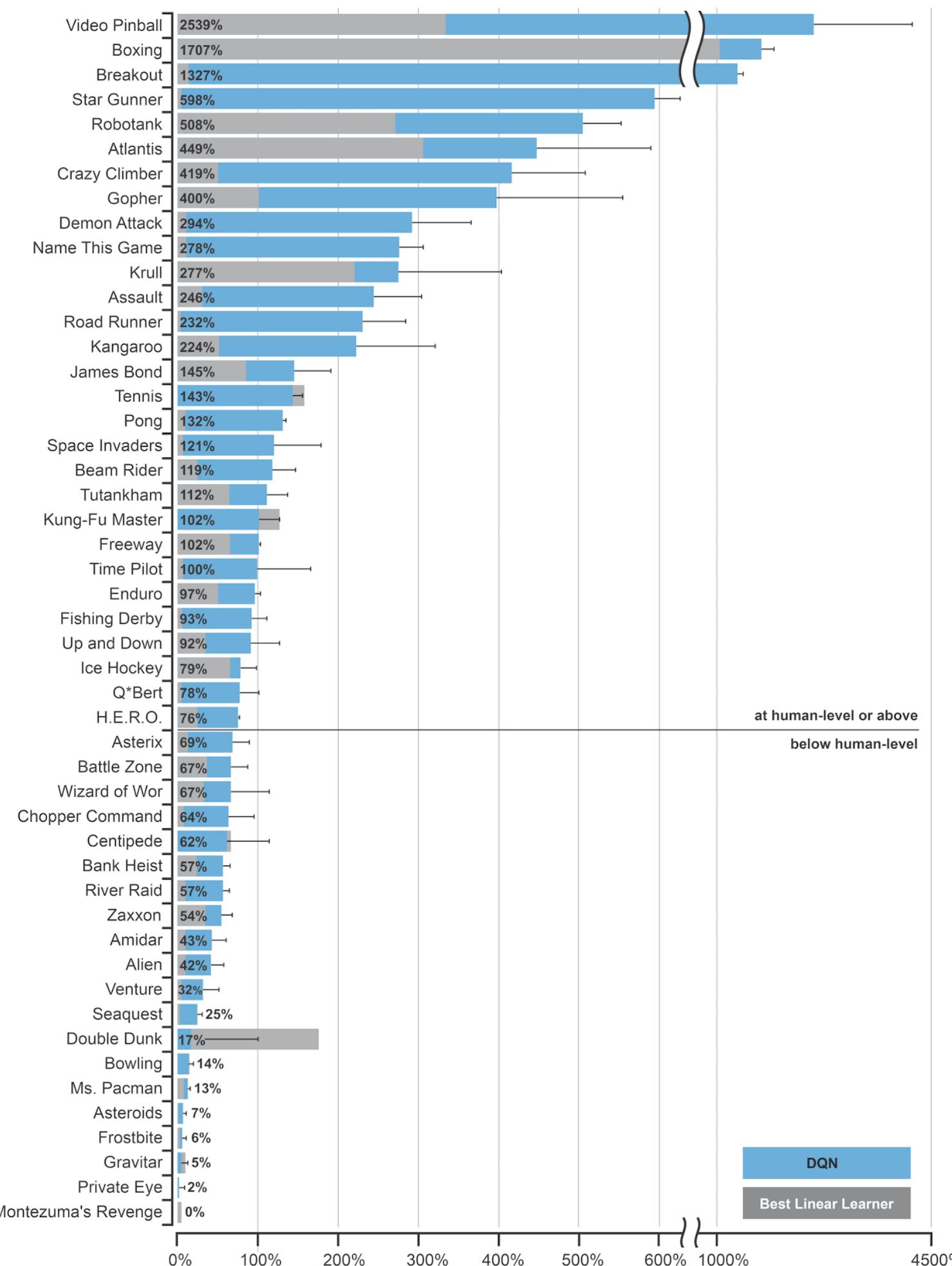
DQN to solve multiple Atari games



DQN to solve multiple Atari games



DQN to solve multiple Atari games



- The DQN network was trained to solve 49 different Atari 2600 games **with the same architecture and hyperparameters**.
- In most of the games, the network reaches **super-human** performance.
- Some games are still badly performed (e.g. Montezuma's revenge), as they require long-term planning.
- It was the first RL algorithm able to learn different tasks (no free lunch theorem).
- The 2015 paper in Nature started the hype for deep RL.

2 - Double DQN

Deep Reinforcement Learning with Double Q-learning

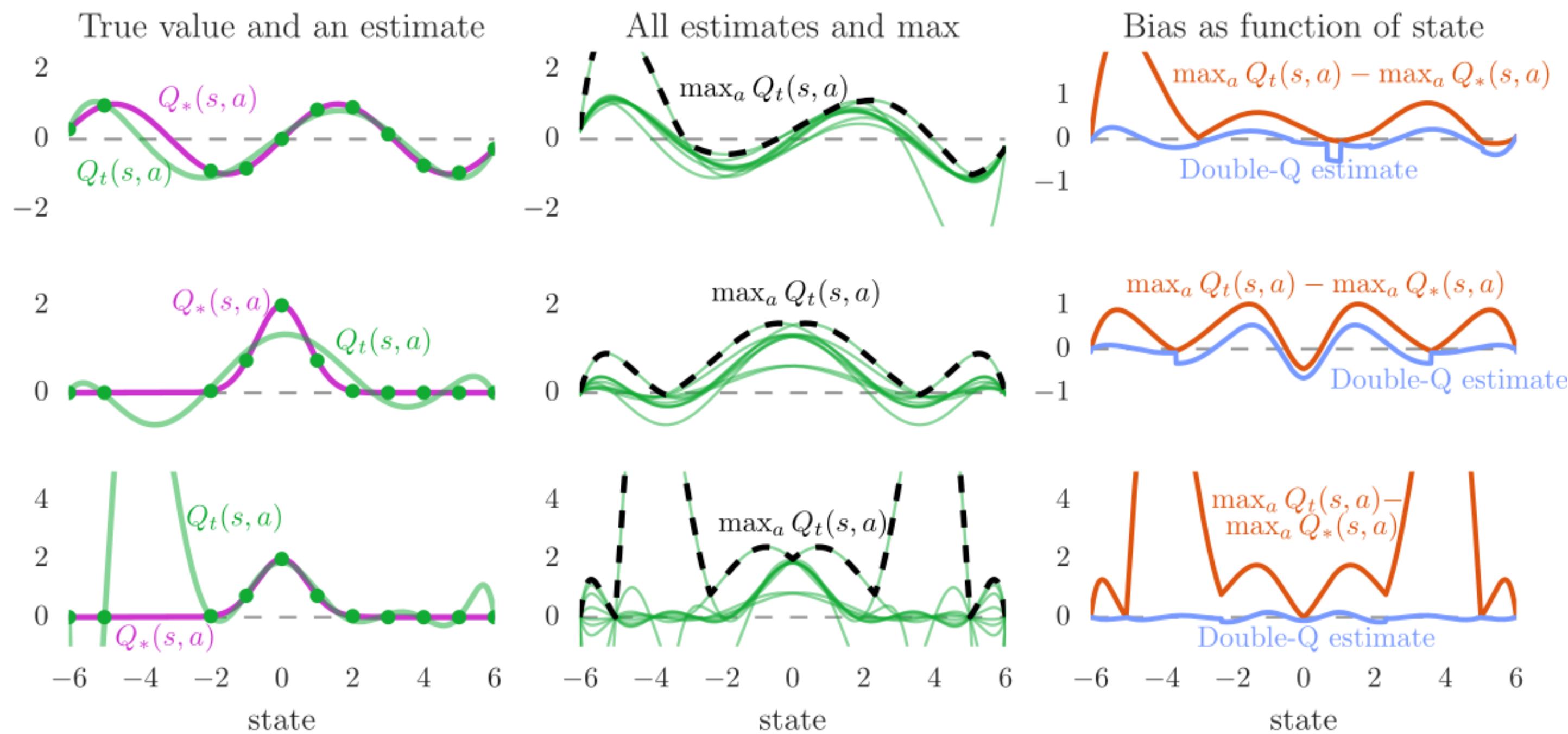
Hado van Hasselt and Arthur Guez and David Silver
Google DeepMind

Double DQN

- Q-learning methods, including DQN, tend to **overestimate** Q-values, especially for the non-greedy actions:

$$Q_\theta(s, a) > Q^\pi(s, a)$$

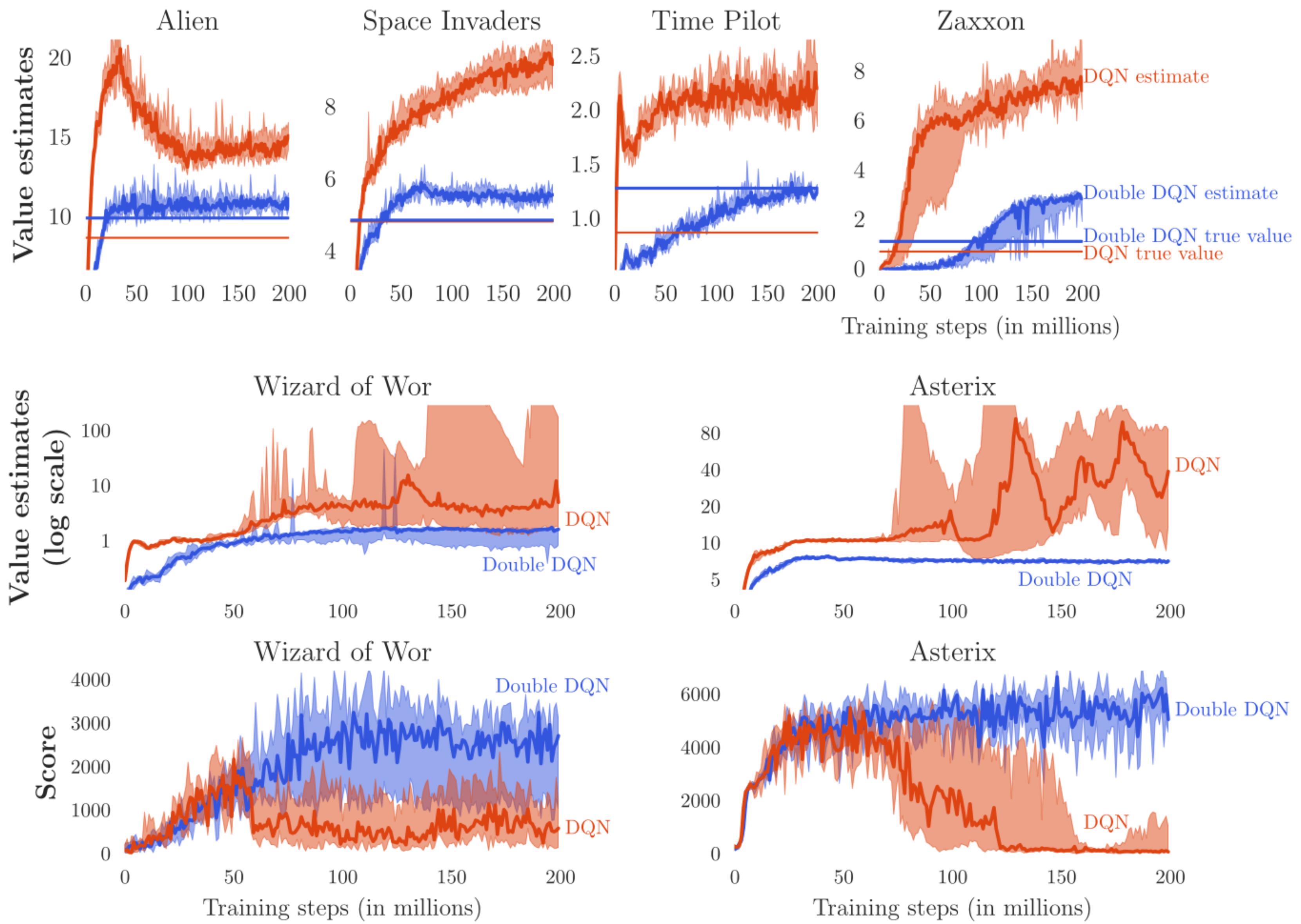
- This does not matter much in action selection, as we apply ϵ -greedy or softmax on the Q-values anyway, but it may make learning slower (sample complexity) and less optimal.



Double DQN

- To avoid optimistic estimations, the target is computed by both the value network θ and the target network θ' :
 - **Action selection:** The next greedy action a^* is calculated by the **value network** θ (current policy):
$$a^* = \operatorname{argmax}_{a'} Q_\theta(s', a')$$
 - **Action evaluation:** Its Q-value for the target is calculated using the **target network** θ' (older values):
$$t = r + \gamma Q_{\theta'}(s', a^*)$$
- This gives the following loss function for **double DQN** (DDQN):

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathcal{D}}[(r + \gamma Q_{\theta'}(s', \operatorname{argmax}_{a'} Q_\theta(s', a'))) - Q_\theta(s, a))^2]$$



3 - Prioritized Experience Replay

Published as a conference paper at ICLR 2016

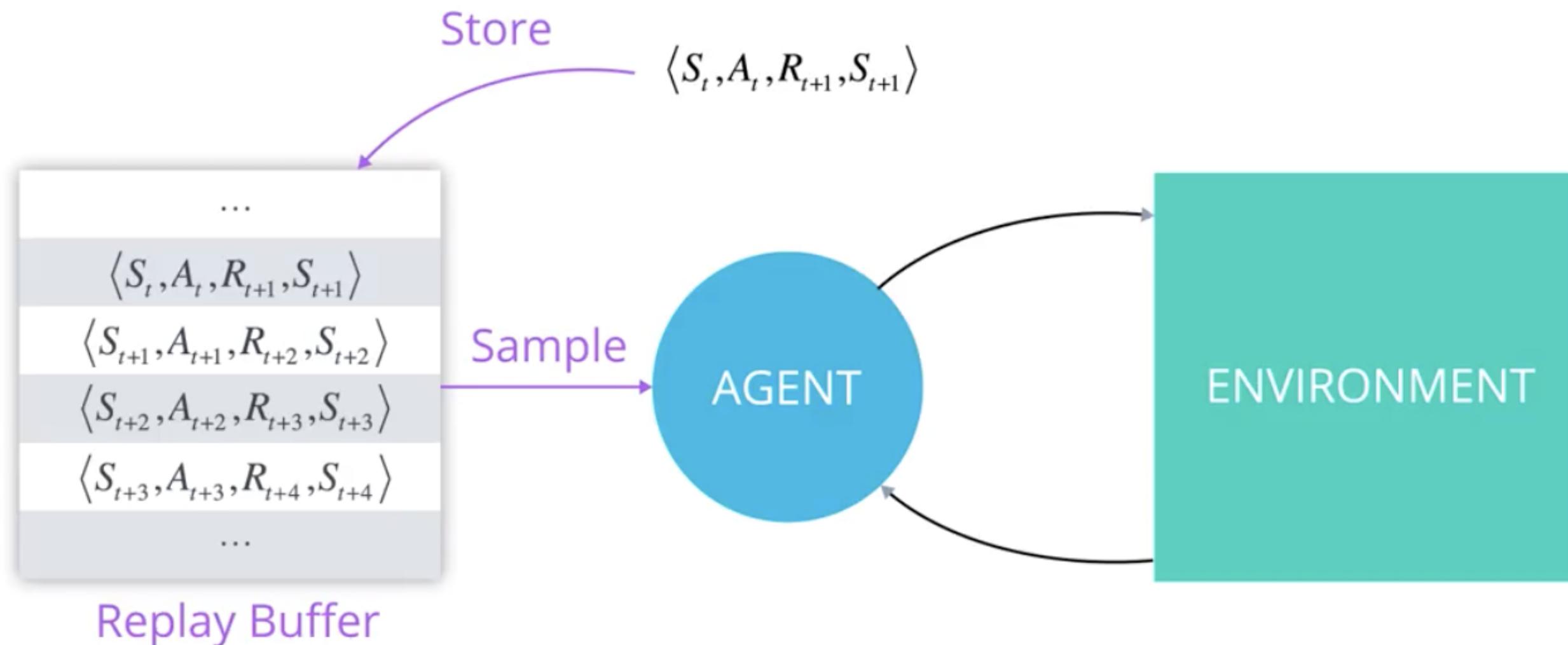
PRIORITIZED EXPERIENCE REPLAY

Tom Schaul, John Quan, Ioannis Antonoglou and David Silver

Google DeepMind

{schaul, johnquan, ioannisa, davidsilver}@google.com

Prioritized Experience Replay

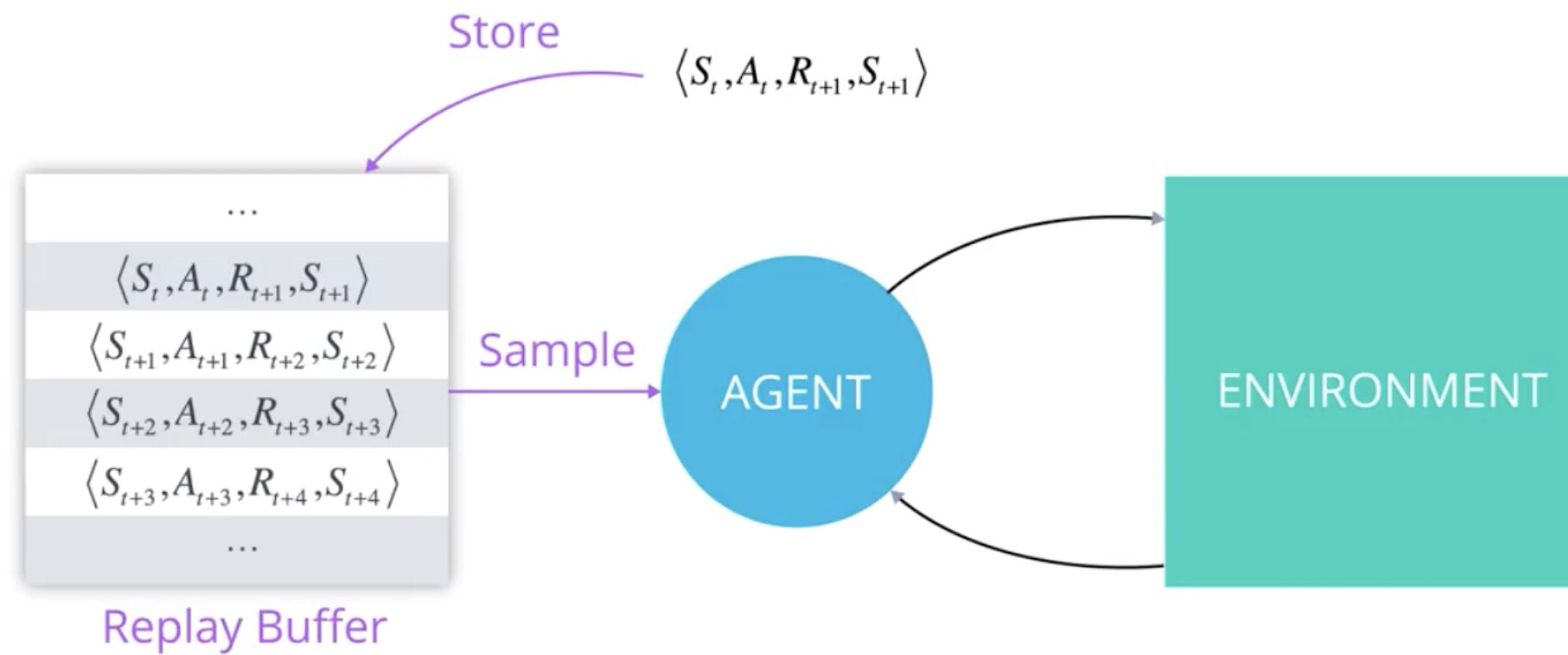


Source: https://parksurk.github.io/deep/reinforcement/learning/drlInd_2-4_value_based_methods-post/

- The **experience replay memory** or **replay buffer** is used to store the last 1M transitions (s, a, r, s') .
- The learning algorithm **randomly samples** a minibatch of size K to update its parameters.
- Not all transitions are interesting:
 - Some transitions were generated by a very old policy, the current policy won't take them anymore.
 - Some transitions are already well predicted: the TD error is small, there is nothing to learn from.

$$\delta_t = r_{t+1} + \gamma \max_{a'} Q_\theta(s_{t+1}, a_{t+1}) - Q_\theta(s_t, a_t) \approx 0$$

Prioritized Experience Replay



Source: https://parksurk.github.io/deep/reinforcement/learning/drlInd_2-4_value_based_methods-post/

- The experience replay memory makes learning very **slow**: we need a lot of samples to learn something useful:
 - High **sample complexity**.
- We need a smart mechanism to preferentially pick the transitions that will boost learning the most, without introducing a bias.
 - **Prioritized sweeping** is actually a quite old idea:

Moore and Atkeson (1993). Prioritized sweeping: Reinforcement learning with less data and less time. Machine Learning, 13(1):103–130.

Prioritized Experience Replay

- The idea of **prioritized experience replay** (PER) is to sample in priority those transitions whose TD error is the highest:

$$\delta_t = r_{t+1} + \gamma \max_{a'} Q_\theta(s_{t+1}, a_{t+1}) - Q_\theta(s_t, a_t)$$

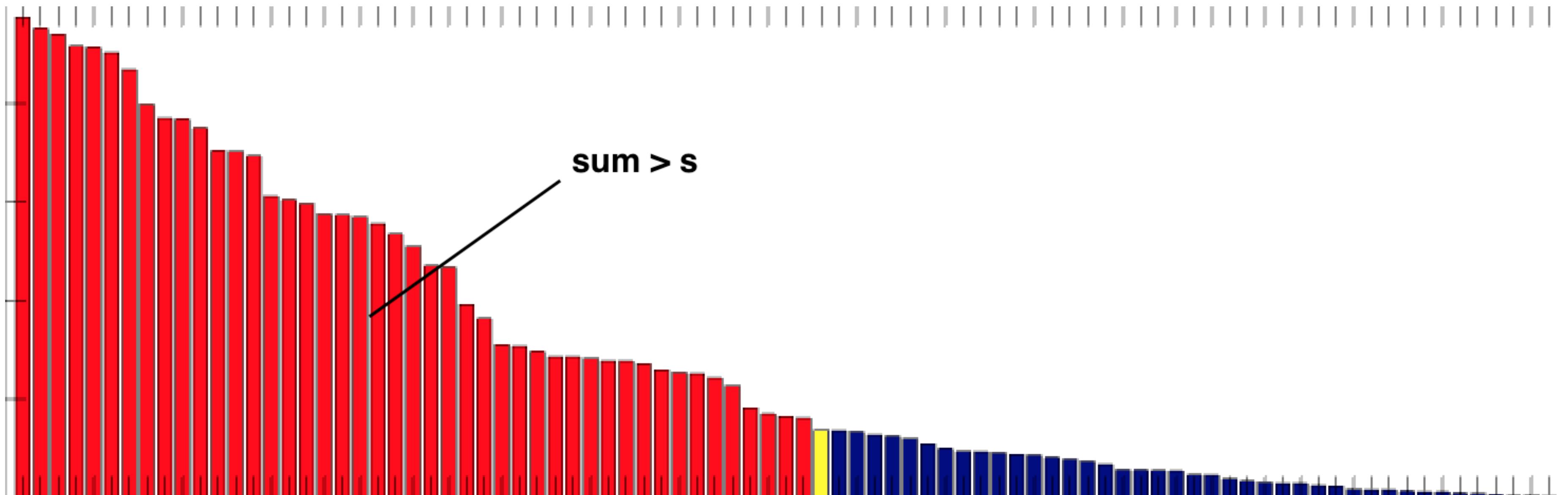
- In practice, we insert the transition (s, a, r, s', δ) into the replay buffer.
- To create a minibatch, the sampling algorithm select a transition k based on the probability:

$$P(k) = \frac{(|\delta_k| + \epsilon)^\alpha}{\sum_k (|\delta_k| + \epsilon)^\alpha}$$

- ϵ is a small parameter ensuring that transition with no TD error still get sampled from time to time.
- α allows to change the behavior from uniform sampling ($\alpha = 0$, as in DQN) to fully prioritized sampling ($\alpha = 1$). α should be annealed from 0 to 1 during training.
- Think of it as a “kind of” **softmax** over the TD errors.
- After the samples have been used for learning, their TD error δ is updated in the PER.

Prioritized Experience Replay

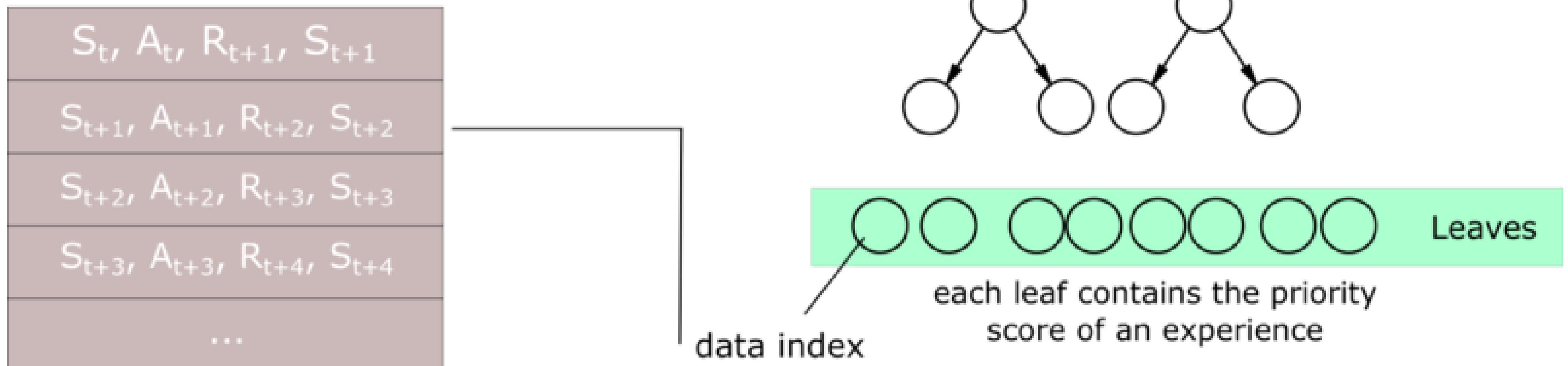
- The main drawback is that inserting and sampling can be computationally expensive if we simply sort the transitions based on $(|\delta_k| + \epsilon)^\alpha$:
 - Insertion: $\mathcal{O}(N \log N)$.
 - Sampling: $\mathcal{O}(N)$.



Source: <https://jaromiru.com/2016/11/07/lets-make-a-dqn-double-learning-and-prioritized-experience-replay/>

Prioritized Experience Replay

- Using binary **sumtrees**, prioritized experience replay can be made efficient in both insertion ($\mathcal{O}(\log N)$) and sampling ($\mathcal{O}(1)$).
- Instead of a linear queue, we use a binary tree to store the transitions.
- Details in a real computer science course...

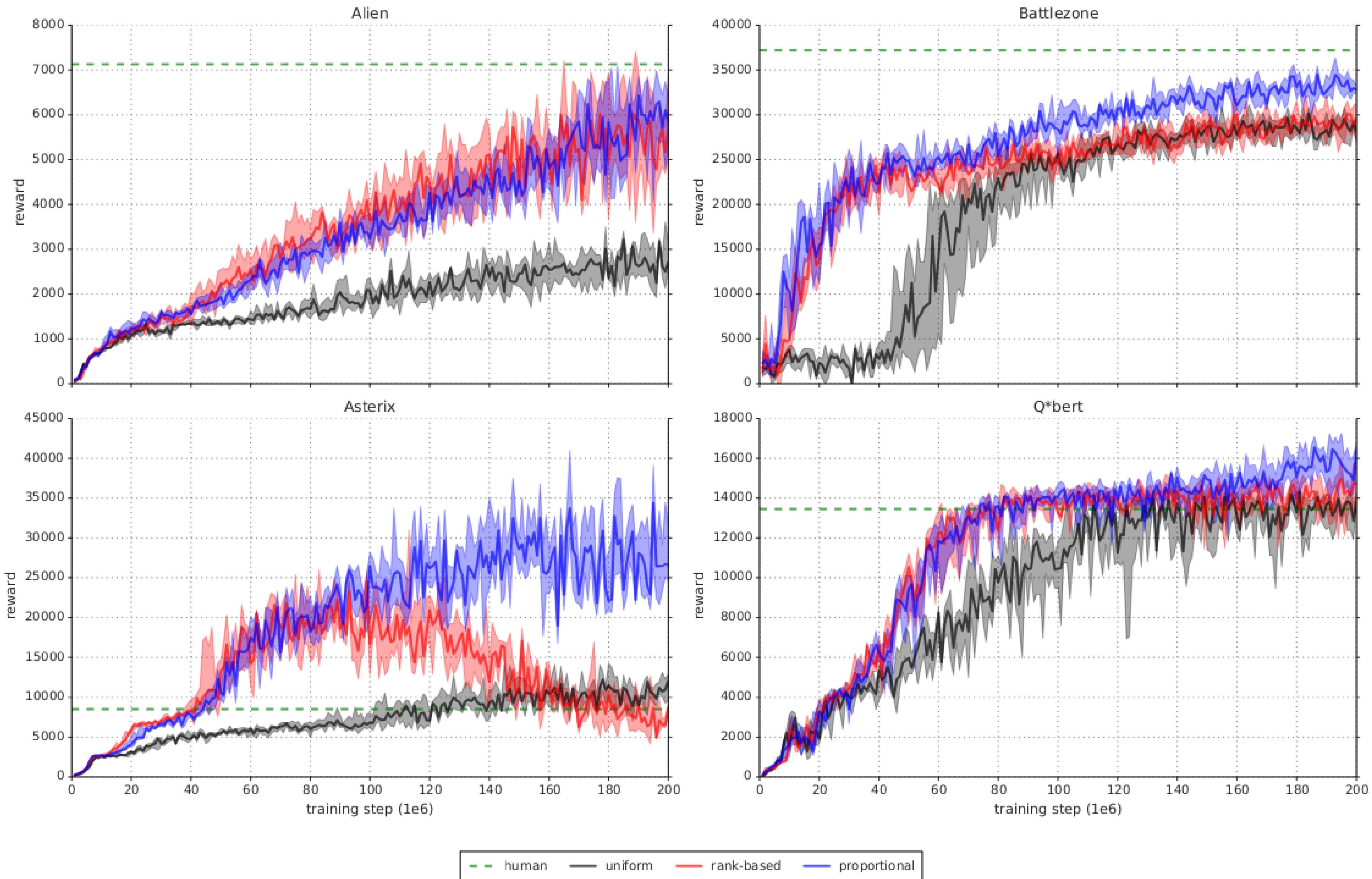


Data

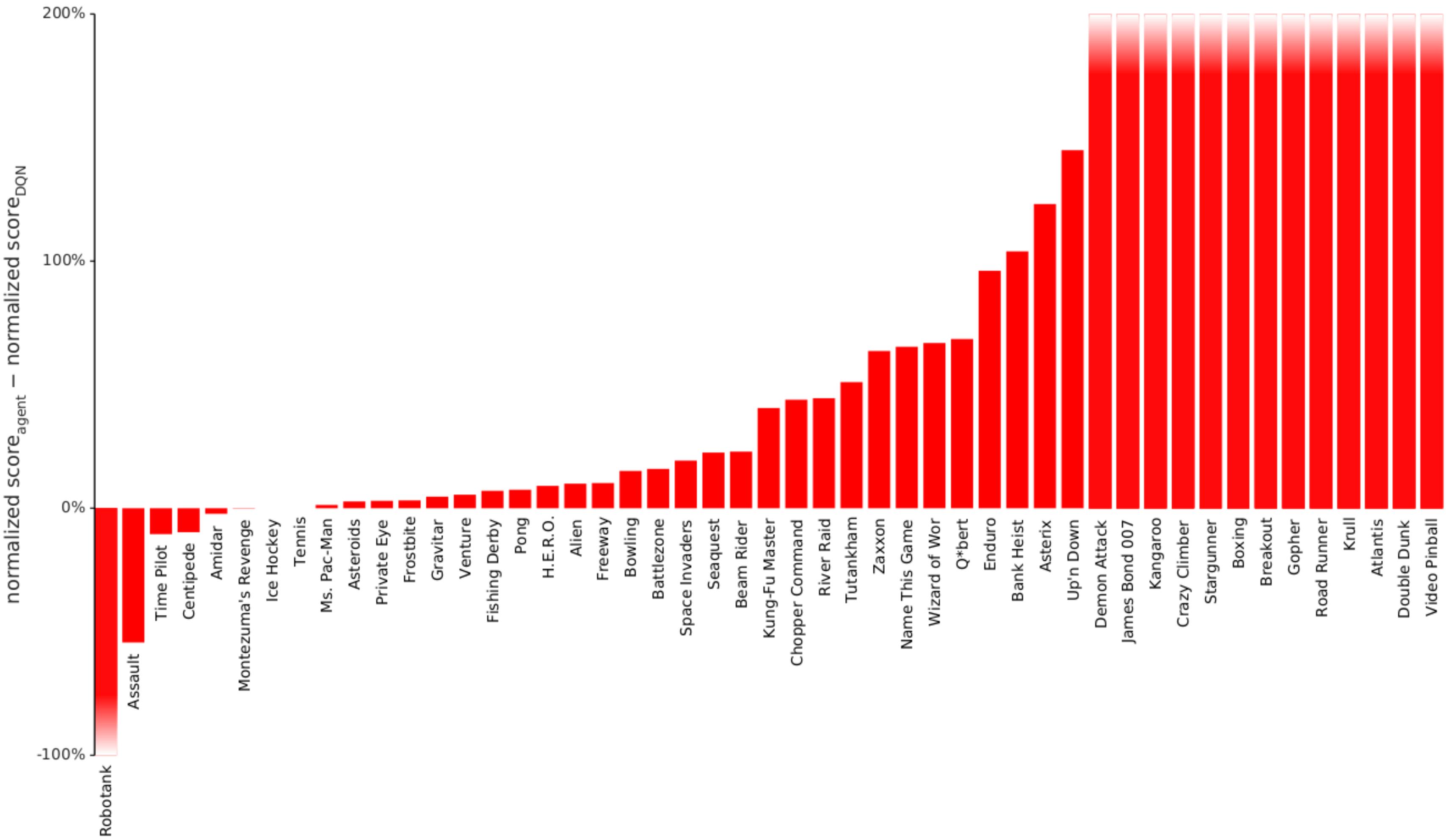
SumTree

Source: <https://www.freecodecamp.org/news/improvements-in-deep-q-learning-dueling-double-dqn-prioritized-experience-replay-and-fixed-58b130cc5682/>

Prioritized Experience Replay



Prioritized Experience Replay



4 - Dueling networks

Dueling Network Architectures for Deep Reinforcement Learning

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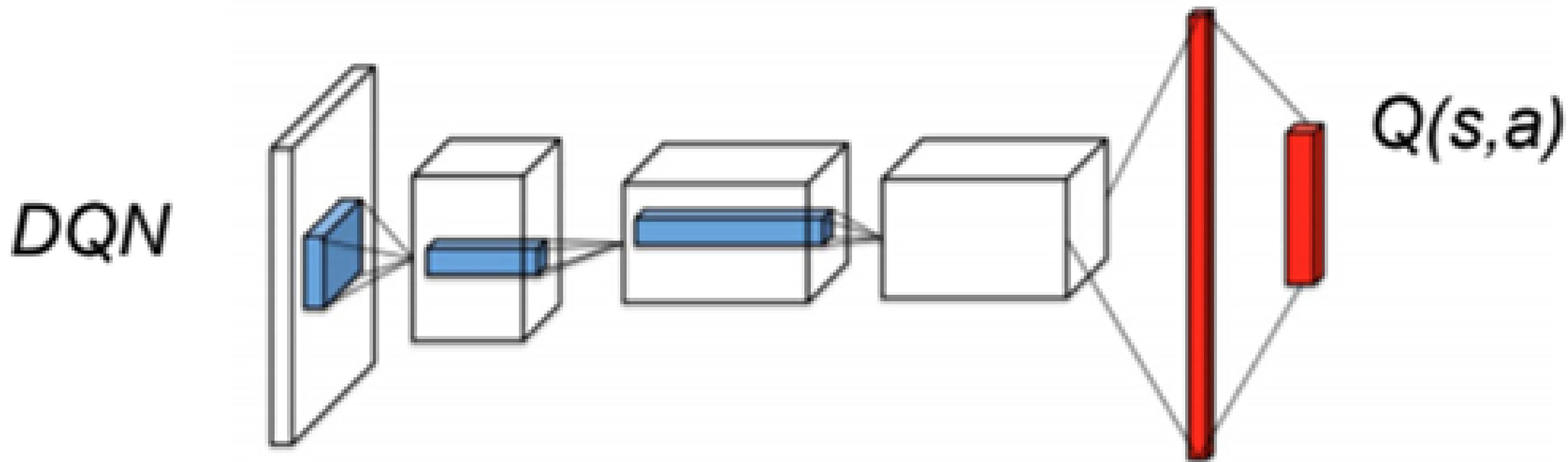
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Dueling networks

- DQN and its variants learn to predict directly the Q-value of each available action.



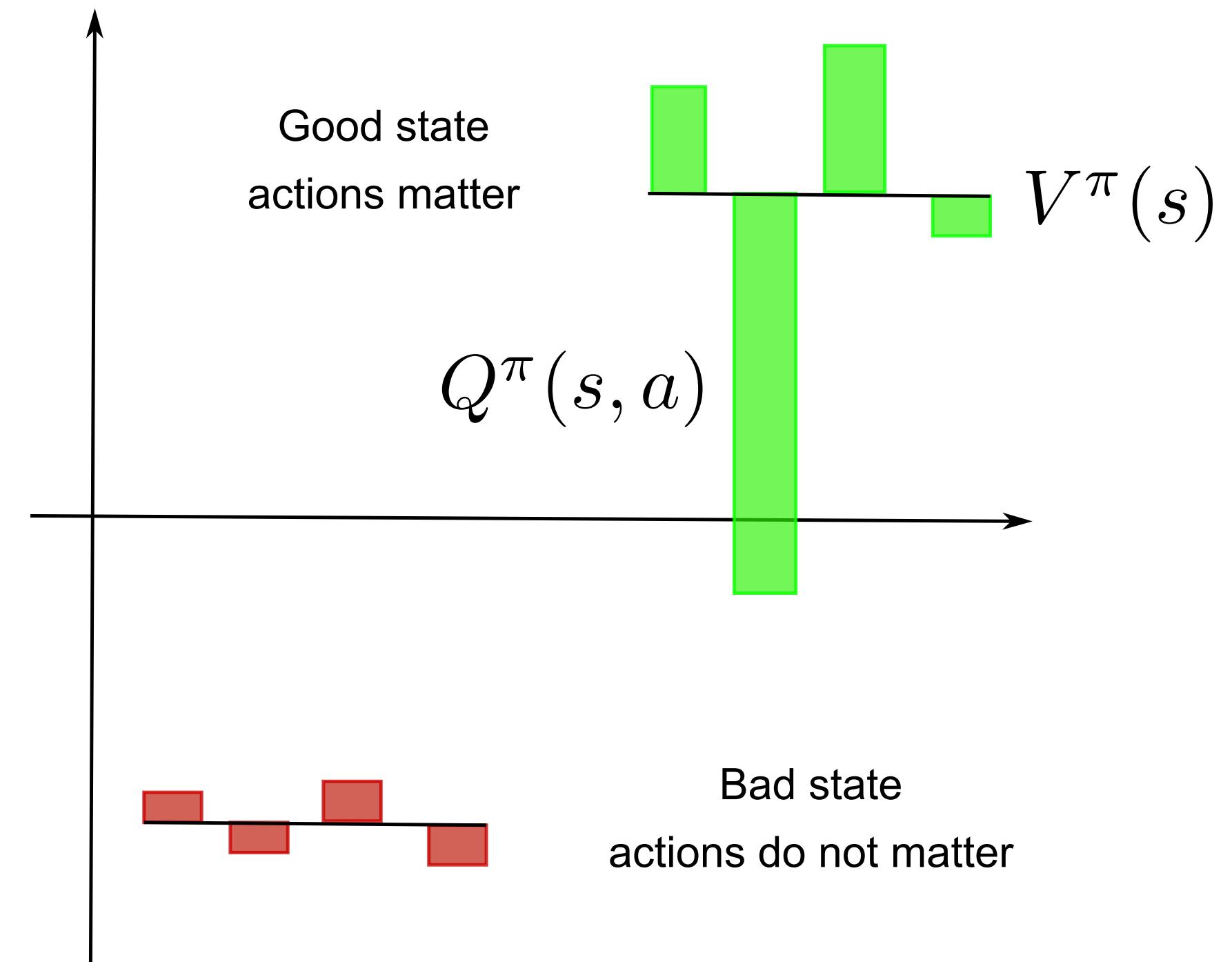
- Several problems with predicting Q-values with a DNN:
 - The Q-values can take high values, especially with different values of γ .
 - The Q-values have a high variance, between the minimum and maximum returns obtained during training.
 - For a transition (s_t, a_t, s_{t+1}) , a single Q-value is updated, not all actions in s_t .

Dueling networks

- Enduro game.

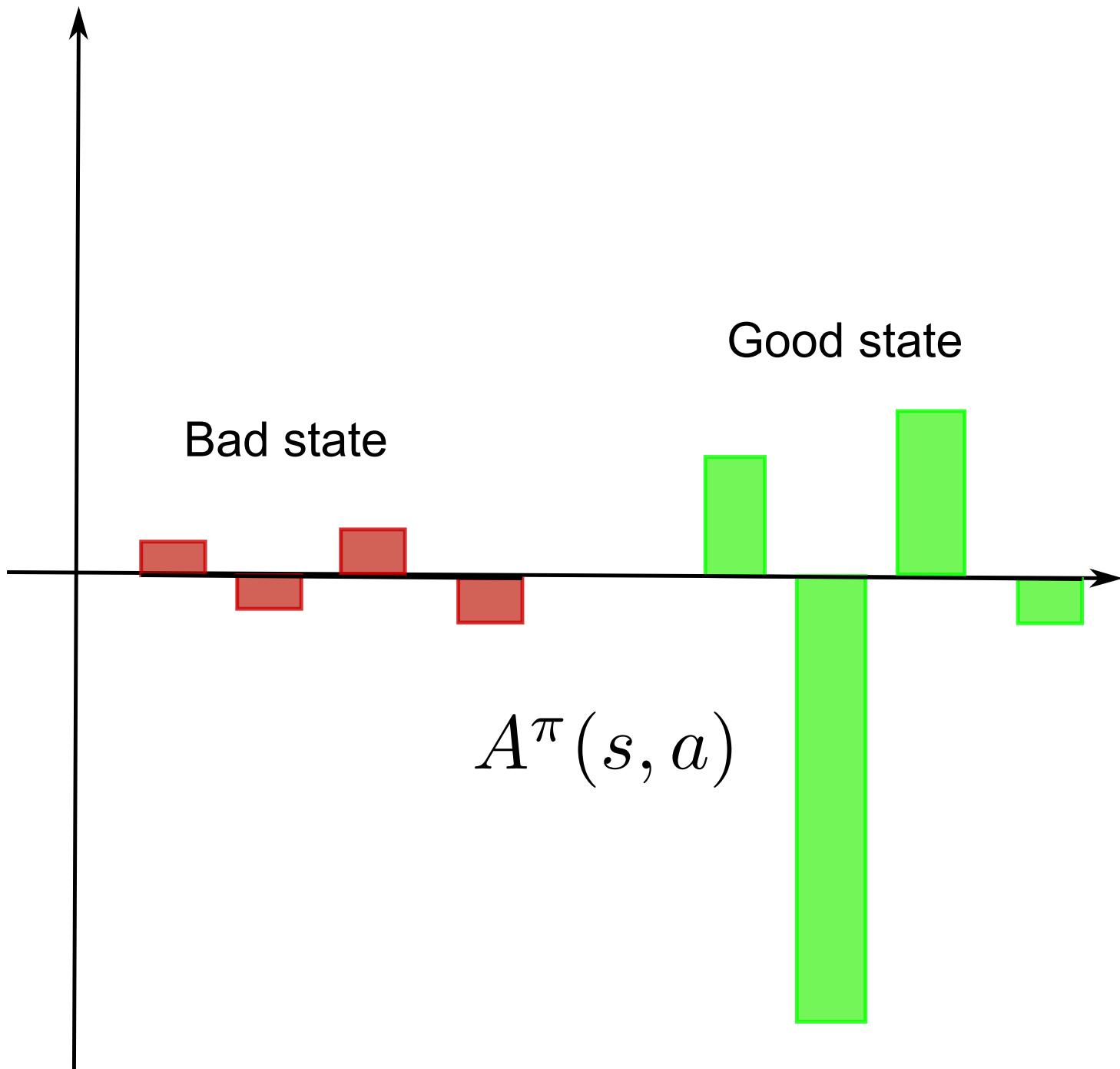


Source: <https://gfycat.com/clumsypaleimpala>



- The exact Q-values of all actions are not equally important.
 - In **bad** states (low $V^\pi(s)$), you can do whatever you want, you will lose.
 - In neutral states, you can do whatever you want, nothing happens.
 - In **good** states (high $V^\pi(s)$), you need to select the right action to get rewards, otherwise you lose.

Advantage functions



- An important notion is the **advantage** $A^\pi(s, a)$ of an action:

$$A^\pi(s, a) = Q^\pi(s, a) - V^\pi(s)$$

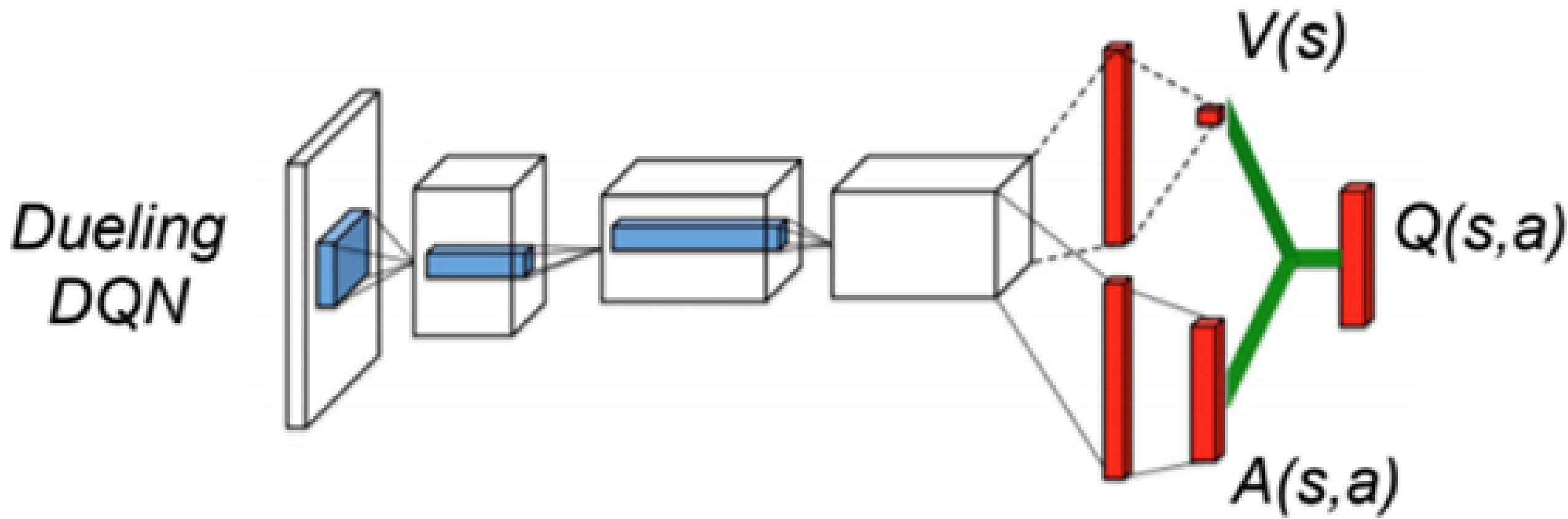
- It tells how much return can be expected by taking the action a in the state s , **compared** to what is usually obtained in s with the current policy.
- If a policy π is deterministic and always selects a^* in s , we have:

$$A^\pi(s, a^*) = 0$$

$$A^\pi(s, a \neq a^*) < 0$$

- This is particularly true for the optimal policy.
- But if we have separate estimates $V_\varphi(s)$ and $Q_\theta(s, a)$, some actions may have a positive advantage.
- Advantages have **less variance** than Q-values.

Dueling networks



- In **dueling networks**, the network is forced to decompose the estimated Q-value $Q_\theta(s, a)$ into a state value $V_\alpha(s)$ and an advantage function $A_\beta(s, a)$:

$$Q_\theta(s, a) = V_\alpha(s) + A_\beta(s, a)$$

- The parameters α and β are just two shared subparts of the NN θ .
- The loss function

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathcal{D}}[(r + \gamma Q_{\theta'}(s', \text{argmax}_{a'} Q_\theta(s', a')) - Q_\theta(s, a))^2]$$

is exactly the same as in (D)DQN: only the internal structure of the NN changes.

Unidentifiability

- The Q-values are the sum of two functions: $Q_\theta(s, a) = V_\alpha(s) + A_\beta(s, a)$
- However, the sum is **unidentifiable**:

$$\begin{aligned} Q_\theta(s, a) &= 10 = 1 + 9 \\ &= 2 + 8 \\ &= 3 + 7 \end{aligned}$$

- To constrain the sum, (Wang et al. 2016) propose that the greedy action w.r.t the advantages should have an advantage of 0:

$$Q_\theta(s, a) = V_\alpha(s) + (A_\beta(s, a) - \max_{a'} A_\beta(s, a'))$$

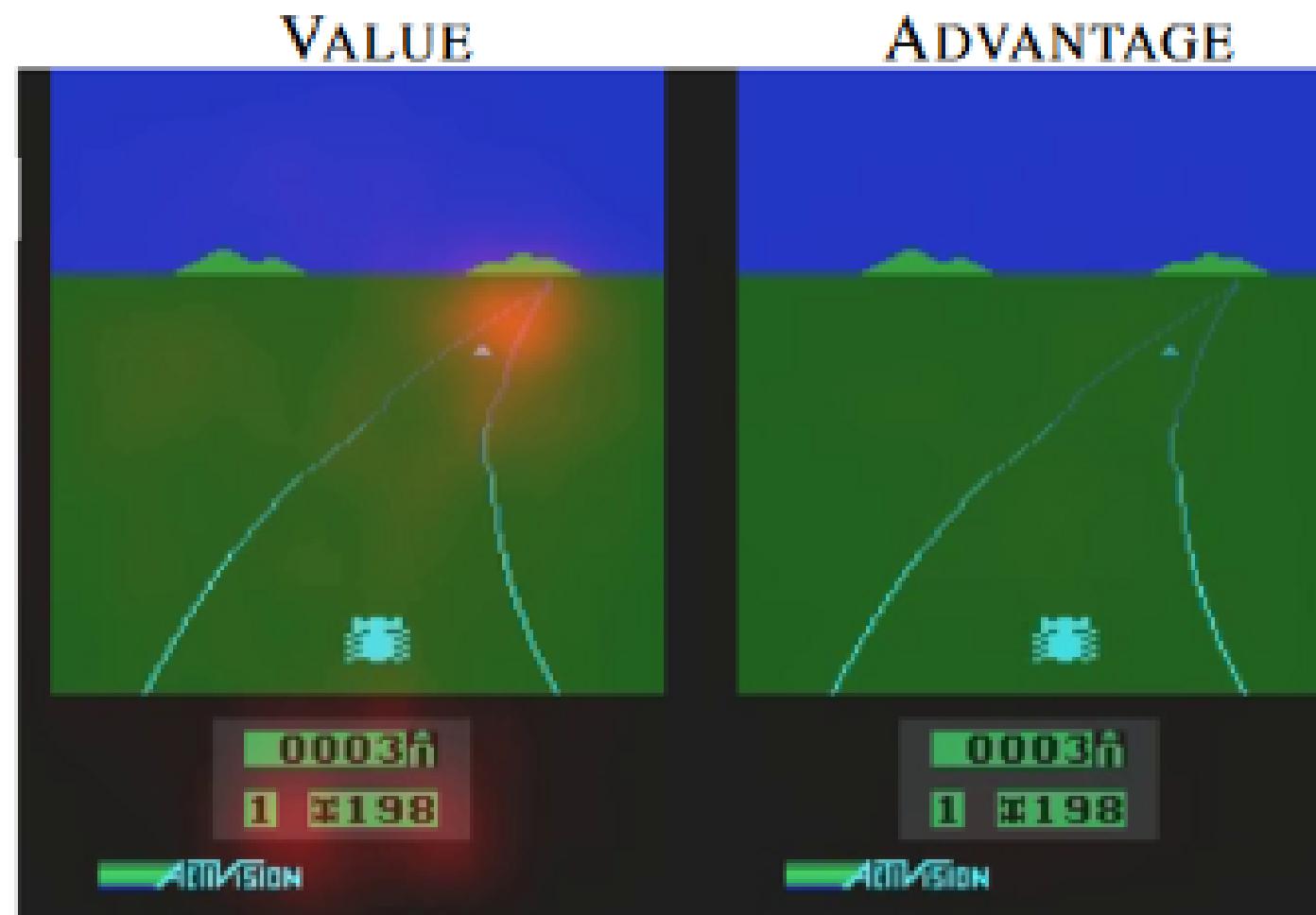
- This way, there is only one solution to the addition. The operation is differentiable, so backpropagation will work.
- (Wang et al. 2016) show that subtracting the mean advantage works better in practice:

$$Q_\theta(s, a) = V_\alpha(s) + (A_\beta(s, a) - \frac{1}{|\mathcal{A}|} \sum_{a'} A_\beta(s, a'))$$

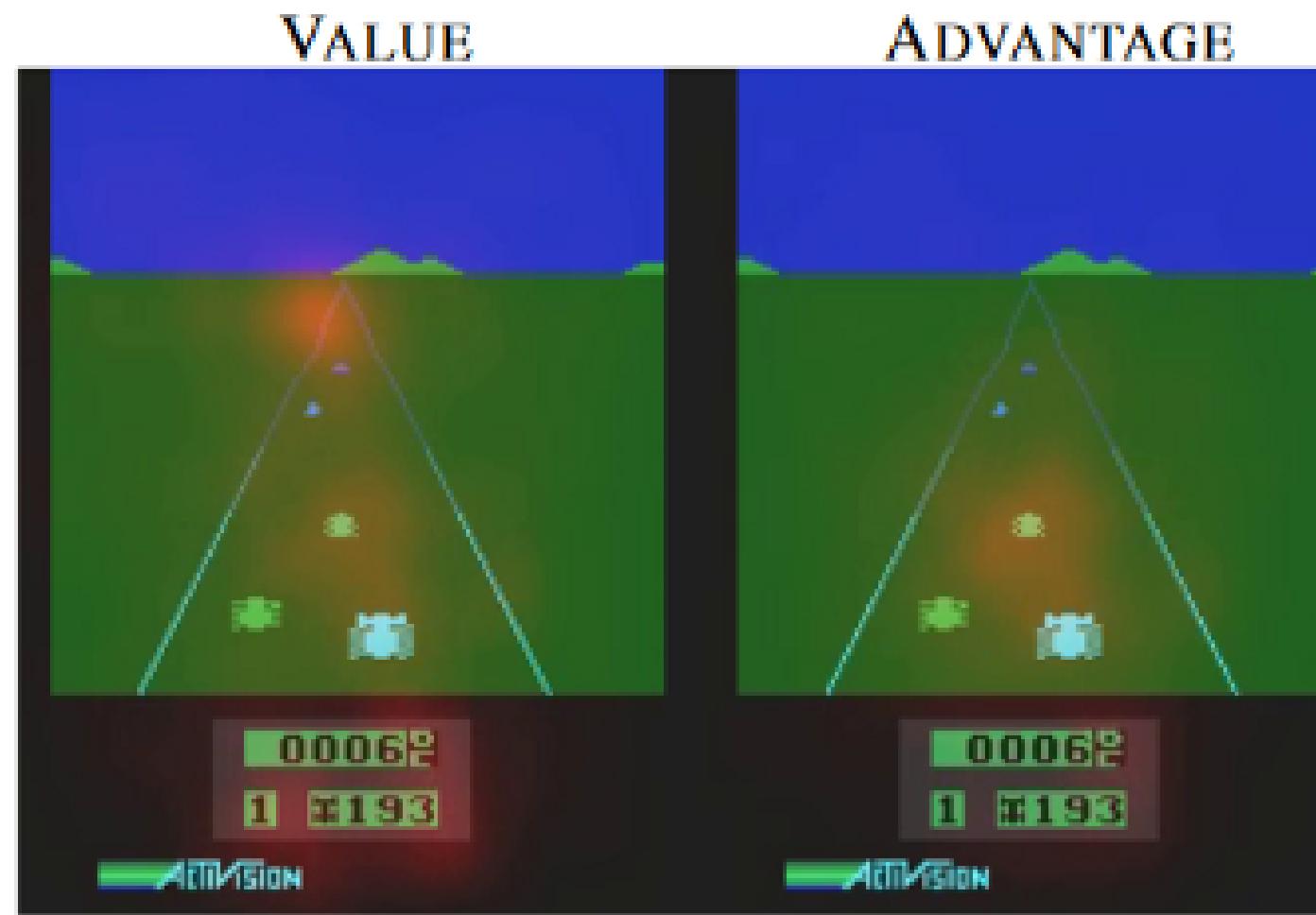
Visualization of the value and advantage functions

- Which pixels change the most the value and advantage functions?

Focus on 2 things:
- The horizon where new cars appear
- On the score

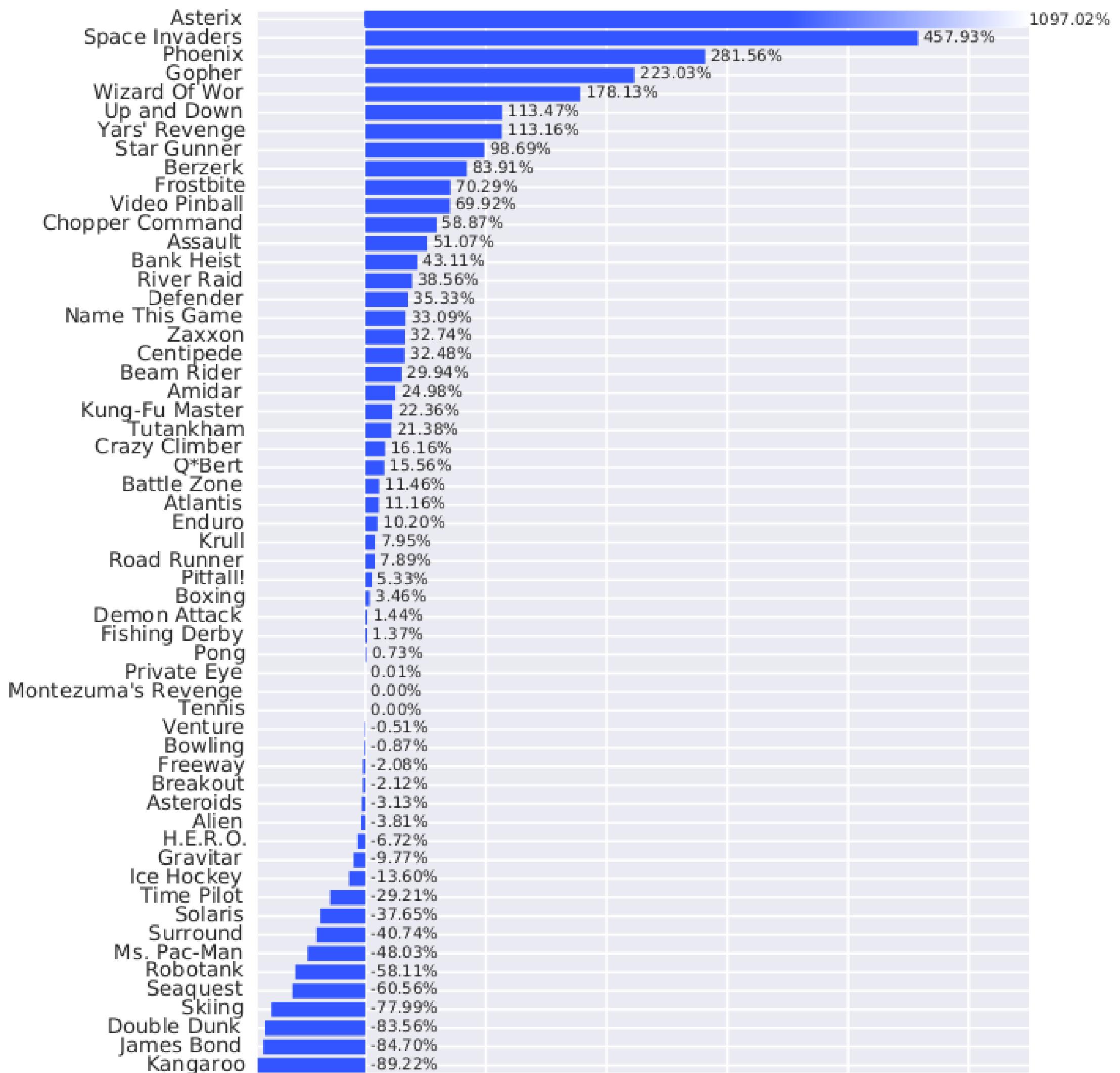


No car in front,
does not pay
much attention
because action
choice making is
not relevant

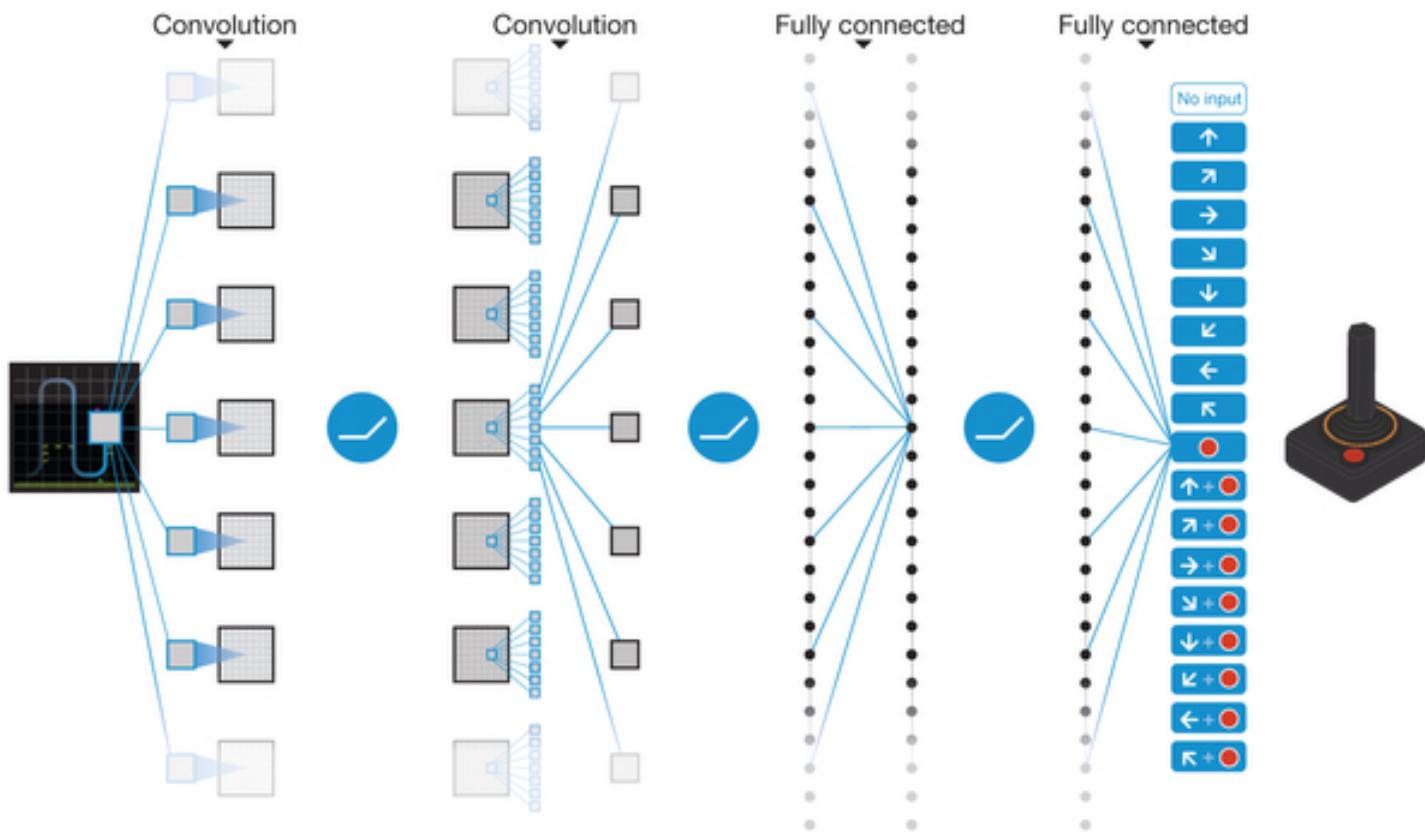


Pays attention to
the front car, in this
case choice
making is crucial
to survive

Improvement over prioritized DDQN



Summary of DQN



- DQN and its early variants (double duelling DQN with PER) are an example of **value-based deep RL**.
- The value $Q_\theta(s, a)$ of each possible action in a given state is approximated by a convolutional neural network.
- The NN has to minimize the mse between the predicted Q-values and the target value corresponding to the Bellman equation:

$$\mathcal{L}(\theta) = \mathbb{E}_{\mathcal{D}}[(r + \gamma Q_{\theta'}(s', \text{argmax}_{a'} Q_\theta(s', a')) - Q_\theta(s, a))^2]$$

- The use of an **experience replay memory** and of **target networks** allows to stabilize learning and avoid suboptimal policies.
- The main drawback of DQN is **sample complexity**: it needs huge amounts of experienced transitions to find a correct policy. The sample complexity come from the deep network itself (gradient descent is iterative and slow), but also from the ERM: it contains 1M transitions, most of which are outdated.
- Only works for **small and discrete action spaces** (one output neuron per action).