# Winter School 2024 Reinforcement Learning

### Advanced Reinforcement Learning

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### Outlook

- Value Function Approximation
- Policy Gradient
- Deep Reinforcement Learning

# Large-Scale Reinforcement Learning

-Reinforcement learning can be used to solve large problems, e.g.

- Backgammon: 10<sup>20</sup> states
- Go: 10<sup>170</sup>
- Robots: continuous state space
- -How can we scale up e.g., the model-free methods for prediction and control seen before?

# Value Function Approximation

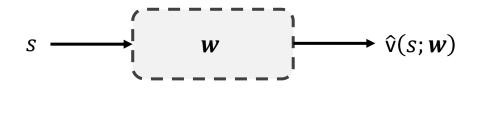
-So far, we have represented value function by a lookup table

- Every state s has an entry V(s)
- Or every state-action pair s, a has an entry Q(s, a)
- Problem with large MDPs:
  - There are too many states and/or actions to store in memory
  - It is too slow to learn the value of each state individually
- -Solution for large MDPs:
  - Estimate value function with function approximation

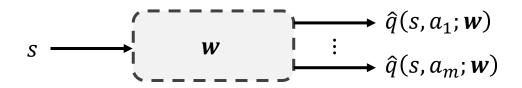
 $\hat{\forall}(s; \mathbf{w}) \approx \lor_{\pi}(s)$  $\hat{q}(s, a; \mathbf{w}) \approx q_{\pi}(s, a)$ 

- Generalize from seen states to unseen states
- Update parameter w using MC or TD learning

### **Types of Value Function Approximation**







### Value Function Approx. By Stochastic Gradient Descent

- Goal: find parameter vector w minimizing mean-squared error between approximate value function  $\hat{v}(s; w)$  and true value function  $v_{\pi}(s)$ 

$$J(\boldsymbol{w}) = \mathbb{E}_{\pi}[(\vee_{\pi}(s) - \hat{\vee}(s; \boldsymbol{w}))^2]$$

- Gradient descent finds a local minimum

$$\Delta \boldsymbol{w} = -\frac{1}{2} a \nabla_{\boldsymbol{w}} J(\boldsymbol{w})$$
$$= \alpha \mathbb{E}_{\pi} [(\vee_{\pi}(s) - \hat{\vee}(s; \boldsymbol{w})) \nabla_{\boldsymbol{w}} \hat{\vee}(s; \boldsymbol{w})]$$

- Stochastic gradient descent samples the gradient

$$\Delta \boldsymbol{w} = \alpha(\vee_{\pi}(s) - \hat{\vee}(s; \boldsymbol{w}))\nabla_{\boldsymbol{w}}\hat{\vee}(s; \boldsymbol{w})$$

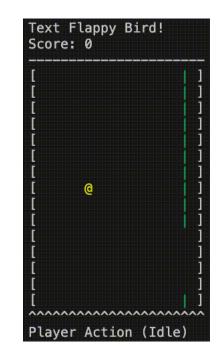
- Expected update is equal to full gradient update

### **Feature Vectors**

-Represent state by a feature vector

$$x(s) = \begin{pmatrix} x_1(s) \\ \vdots \\ x_n(s) \end{pmatrix}$$

- -E.g., Polynomials, Fourier Basis
- -For example:
  - Distance of robot from landmarks
  - Trends in the stock market
  - Piece and pawn configurations in chess



### Linear Value Function Approximation

- Represent value function by a linear combination of features

$$\hat{\mathbf{v}}(s; \mathbf{w}) = x(s)^T \mathbf{w} = \sum_{i=0}^n x_i(s) w_i$$

– Objective function is quadratic in parameters  $\boldsymbol{w}$ 

$$J(w) = \mathbb{E}_{\pi}[(\vee_{\pi}(s) - x(s)^{T} w)^{2}]$$

- Stochastic gradient descent converges on global optimum
- Update rule is particularly simple

$$\nabla_{w} \widehat{\vee}(s; \boldsymbol{w}) = x(s)$$
  
$$\Delta \boldsymbol{w} = \alpha \big( \vee_{\pi}(s) - \widehat{\vee}(s; \boldsymbol{w}) \big) x(s)$$

– Update = step-size × prediction error × feature value

### **Incremental Prediction Algorithms**

- -Assumed true value  $v_{\pi}(s)$  function given by supervisor
- -But in RL there is no supervisor, only rewards
- -In practice, we substitute a target for  $v_{\pi}(s)$

– For MC, the target is the return  $G_t$ 

$$\Delta \boldsymbol{w} = \alpha \big( \boldsymbol{G}_{\boldsymbol{t}} - \hat{\vee}(\boldsymbol{S}_{\boldsymbol{t}}; \boldsymbol{w}) \big) \nabla_{\boldsymbol{w}} \hat{\vee}(\boldsymbol{S}_{\boldsymbol{t}}; \boldsymbol{w})$$

- For TD(0), the target is the TD target  $G_{t:t+1} = R_{t+1} + \gamma \hat{v}(s_{t+1}; w)$ 

$$\Delta \boldsymbol{w} = \alpha \big( \boldsymbol{R}_{t+1} + \gamma \,\hat{\boldsymbol{v}}(\boldsymbol{S}_{t+1}; \boldsymbol{w}) - \hat{\boldsymbol{v}}(\boldsymbol{S}_t; \boldsymbol{w}) \big) \nabla_{\boldsymbol{w}} \hat{\boldsymbol{v}}(\boldsymbol{S}_t; \boldsymbol{w})$$

### Monte-Carlo with Value Function Approximation (1/2)

-Return  $G_t$  is an unbiased, noisy sample of true value  $v_{\pi}(s_t)$ -Can therefore apply supervised learning to "training data":

 $< S_1, G_1 >, < S_2, G_2 >, \dots, < S_T, G_T >$ 

-For example, using linear Monte-Carlo policy evaluation

$$\Delta \boldsymbol{w} = \alpha \big( G_t - \hat{\vee}(S_t; \boldsymbol{w}) \big) \nabla_{\boldsymbol{w}} \hat{\vee}(S_t; \boldsymbol{w}) \\ = \alpha \big( G_t - \hat{\vee}(S_t; \boldsymbol{w}) \big) \boldsymbol{x}(S_t)$$

- -Monte-Carlo evaluation converges to a local optimum
- -Even when using non-linear value function approximation

### Monte-Carlo with Value Function Approximation (2/2)

### Gradient Monte Carlo Algorithm for Estimating $\hat{v} \approx v_{\pi}$

```
Input: the policy \pi to be evaluated
Input: a differentiable function \hat{v} : \mathbb{S} \times \mathbb{R}^d \to \mathbb{R}
Algorithm parameter: step size \alpha > 0
Initialize value-function weights \mathbf{w} \in \mathbb{R}^d arbitrarily (e.g., \mathbf{w} = \mathbf{0})
Loop forever (for each episode):
Generate an episode S_0, A_0, R_1, S_1, A_1, \dots, R_T, S_T using \pi
Loop for each step of episode, t = 0, 1, \dots, T - 1:
\mathbf{w} \leftarrow \mathbf{w} + \alpha [G_t - \hat{v}(S_t, \mathbf{w})] \nabla \hat{v}(S_t, \mathbf{w})
```

### TD Learning with Value Function Approximation (1/2)

- The TD-target  $G_{t:t+1} = R_{t+1} + \gamma \hat{v}(s_{t+1}; w)$  is a biased sample of true value  $v_{\pi}(s_t)$
- Can still apply supervised learning to "training data":

 $< S_1, R_1 + \gamma \hat{v}(S_2; w) >, < S_2, R_3 + \gamma \hat{v}(S_3; w) >, ..., < S_{T-1}, R_T >$ - For example, using linear TD(0)

$$\Delta \boldsymbol{w} = \alpha \big( R + \gamma \,\hat{\vee}(S'; \boldsymbol{w}) - \hat{\vee}(S_t; \boldsymbol{w}) \big) \nabla_{\boldsymbol{w}} \hat{\vee}(S_t; \boldsymbol{w}) = \alpha \delta \boldsymbol{x}(S)$$

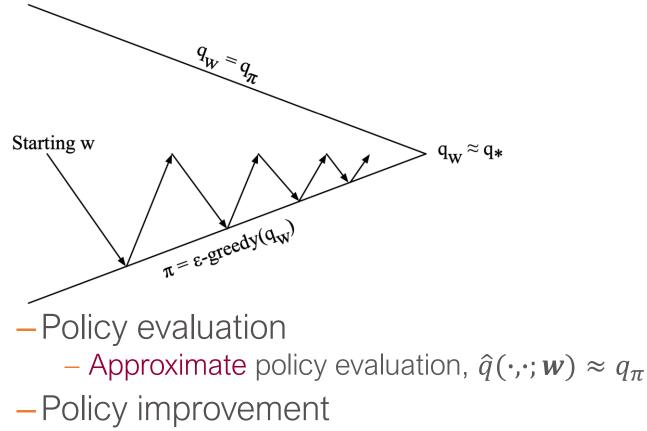
- (Semi-Gradient) We do not consider the effect of changing w on the target - Linear TD(0) converges (close) to global optimum

### TD Learning with Value Function Approximation (2/2)

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

```
Input: the policy \pi to be evaluated
Input: a differentiable function \hat{v}: S^+ \times \mathbb{R}^d \to \mathbb{R} such that \hat{v}(\text{terminal}, \cdot) = 0
Algorithm parameter: step size \alpha > 0
Initialize value-function weights \mathbf{w} \in \mathbb{R}^d arbitrarily (e.g., \mathbf{w} = \mathbf{0})
Loop for each episode:
    Initialize S
    Loop for each step of episode:
        Choose A \sim \pi(\cdot | S)
         Take action A, observe R, S'
        \mathbf{w} \leftarrow \mathbf{w} + \alpha [R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})] \nabla \hat{v}(S, \mathbf{w})
        S \leftarrow S'
    until S is terminal
```

### **Control with Value Function Approximation**



<u>ε-Greedy</u> policy improvement

[David Silver, IRL, UCL 2015]

### **Action-Value Function Approximation**

- Approximate the action-value function

 $\hat{q}(S, A; \boldsymbol{w}) \approx q_{\pi}(S, A)$ 

– Minimize mean-squared error between approximate action-value function  $\hat{q}(S, A, w)$  and true action-value function  $q_{\pi}(S, A)$ 

$$J(\boldsymbol{w}) = \mathbb{E}_{\pi}[(q_{\pi}(S,A) - \hat{q}(S,A;\boldsymbol{w}))^2]$$

- Use stochastic gradient descent to find a local minimum

$$-\frac{1}{2}\nabla_{w}J(\boldsymbol{w}) = (q_{\pi}(S,A) - \hat{q}(S,A;\boldsymbol{w}))\nabla_{w}\hat{q}(S,A;\boldsymbol{w})$$
$$\Delta \boldsymbol{w} = \alpha (q_{\pi}(S,A) - \hat{q}(S,A;\boldsymbol{w}))\nabla_{w}\hat{q}(S,A;\boldsymbol{w})$$

### Linear Action-Value Function Approximation

- Represent state and action by a feature vector

$$x(S,A) = \begin{pmatrix} x_1(S,A) \\ \vdots \\ x_n(S,A) \end{pmatrix}$$

- Represent action-value function by linear combination of features

$$\hat{q}(S, A, \boldsymbol{w}) = x(S, A)^T \boldsymbol{w} = \sum_{i=0}^n x_i(S, A) w_i$$

- Stochastic gradient descent update

$$\nabla_{w}\hat{q}(S,A;\boldsymbol{w}) = x(S,A)$$
  
$$\Delta \boldsymbol{w} = \alpha \big(q_{\pi}(S,A) - \hat{q}(S,A;\boldsymbol{w})\big) x(S,A)$$

### Incremental Control Algorithms

-Like prediction, we must substitute a target for  $q_{\pi}(S, A)$ - For MC, the target is the return  $G_t$ 

$$\Delta \boldsymbol{w} = \alpha \big( \boldsymbol{G}_{\boldsymbol{t}} - \hat{q}(S_t, A_t; \boldsymbol{w}) \big) \nabla_{\boldsymbol{w}} \hat{q}(S_t A_t; \boldsymbol{w})$$

- For TD(0), the target is the TD target  $G_{t:t+1} = R_{t+1} + \gamma \hat{v}(s_{t+1}; w)$ 

$$\Delta \boldsymbol{w} = \alpha \big( \boldsymbol{R}_{t+1} + \gamma \, \hat{\vee} (\boldsymbol{S}_{t+1}; \boldsymbol{w}) - \hat{q}(\boldsymbol{S}_t, \boldsymbol{A}_t; \boldsymbol{w}) \big) \nabla_{\boldsymbol{w}} \hat{q}(\boldsymbol{S}_t \boldsymbol{A}_t; \boldsymbol{w})$$

## Policy-Based Reinforcement Learning

 Previously we approximated the value or action-value function using parameters w

> $\vee_w(s) \approx \vee_\pi(s)$  $q_w(s,a) \approx q_\pi(s,a)$

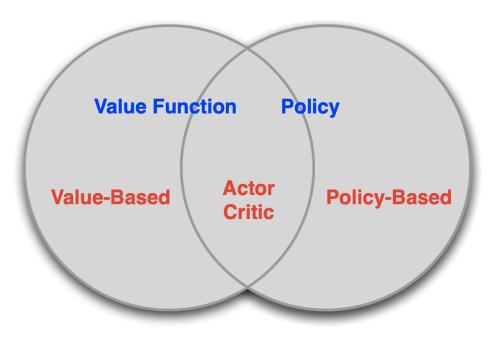
A policy was generated directly from the value function
 – e.g., using ε-greedy

-We now directly parametrize the policy  $\pi_{\theta}(s, a) = \mathbb{P}[a \mid s; \theta]$ 

## Value-Based and Policy-Based RL

### -Value Based

- Learnt Value Function
- Implicit policy
   (e.g., ε-greedy)
- -Policy Based
  - No Value Function
  - Learnt Policy
- -Actor-Critic
  - Learnt Value Function
  - Learnt Policy



# Advantages of Policy-Based RL

### -Advantages:

- Better convergence properties (in contrast to value function approximation that can oscillate in some configurations)
- Effective in high-dimensional or continuous action spaces
- Can learn stochastic policies
- -Disadvantages:
  - Typically converge to a local rather than global optimum
  - Evaluating a policy is typically inefficient and high variance

# **Policy Optimization**

- -Policy based reinforcement learning is an optimization problem
- -Find  $\theta$  that maximizes  $J(\theta)$
- -Many possible optimization approaches
  - Gradient-free (e.g., Hill climbing, Genetic algorithms, etc.)
  - Gradient-based (e.g., Gradient descent)
- -We focus on gradient descent, many extensions possible

# Policy Gradient

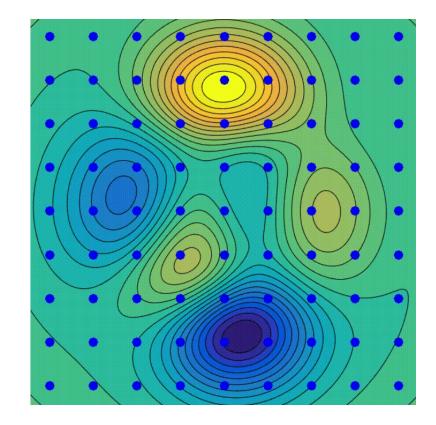
- Let  $J(\theta)$  be any policy objective function
- Policy gradient algorithms search for a local maximum in  $J(\theta)$  by ascending the gradient of the policy, w.r.t. parameters  $\theta$

$$\Delta \theta = a \nabla_{\theta} J(\theta)$$

-Where  $\nabla_{\theta} J(\theta)$  is the policy gradient

$$\nabla_{\theta} J(\boldsymbol{\theta}) = \begin{pmatrix} \frac{\partial J(\theta)}{\partial \theta_1} \\ \vdots \\ \frac{\partial J(\theta)}{\partial \theta_n} \end{pmatrix}$$

– and  $\alpha$  is a step-size parameter



# **Policy Objective Functions**

- Goal: given policy  $\pi_{\theta}(s, a)$  with parameters  $\theta$ , find best  $\theta$
- But how do we measure the quality of a policy  $\pi_{\theta}$ ?
- In episodic environments we can use the start value

$$J_1(\theta) = \mathbf{V}^{\pi_{\theta}}(s_1) = \mathbb{E}_{\pi_{\theta}}[\mathsf{v}_1]$$

- In continuing environments, we can use the average value

$$J_{avV}(\theta) = \sum_{s} d^{\pi_{\theta}}(s) \, \mathrm{V}^{\pi_{\theta}}(s)$$

- Or the average reward per time-step

$$J_{avR}(\theta) = \sum_{s} d^{\pi_{\theta}}(s) \sum_{a} \pi_{\theta}(s, a) \mathcal{R}_{s}^{a}$$

where  $d^{\pi_{\theta}}(s)$  is stationary distribution of Markov chain for  $\pi_{\theta}$ 

## One-Step MDPs

- Consider a simple class of one-step MDPs
  - Starting in state  $s \sim d(s)$
  - Terminating after one time-step with reward  $r = \mathcal{R}_s^a$

- Use likelihood ratios to compute the policy gradient

 $\nabla_{\theta} \pi_{\theta}(s, a) = \pi_{\theta}(s, a) \frac{\nabla_{\theta} \pi_{\theta}(s, a)}{\pi_{\theta}(s, a)}$  $= \pi_{\theta}(s, a) \nabla_{\theta} \log \pi_{\theta}(s, a)$ 

$$J(\theta) = \mathbb{E}_{\pi_{\theta}}[r]$$
  
=  $\sum_{s \in S} d(s) \sum_{a \in A} \pi_{\theta}(s, a) \mathcal{R}_{s}^{a}$   
 $\nabla_{\theta} J(\theta) = \sum_{s \in S} d(s) \sum_{a \in A} \pi_{\theta}(s, a) \nabla_{\theta} \log \pi_{\theta}(s, a) \mathcal{R}_{s}^{a}$   
=  $\mathbb{E}_{\pi_{\theta}}[\nabla_{\theta} \log \pi_{\theta}(s, a) r]$ 

# Policy Gradient Theorem

- The policy gradient theorem generalizes the likelihood ratio approach to multi-step MDPs
- Replaces instantaneous reward r with long-term value  $q_{\pi}(s, a)$
- Policy gradient theorem applies to start state objective, average reward and average value objective

#### Theorem (Policy Gradient)

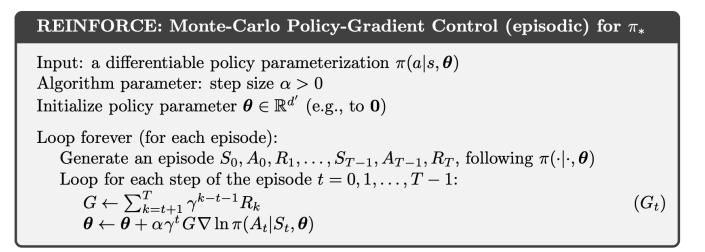
For any differentiable policy  $\pi_{\theta}(s, a)$  and for any of the policy objective functions  $J = J_1, J_{avR}$  or  $\frac{1}{1-v}J_{avV}$  the policy gradient is

 $\nabla_{\theta} J(\theta) = \mathbb{E}_{\pi_{\theta}} [\nabla_{\theta} \log \pi_{\theta}(s, a) q_{\pi}(s, a)]$ 

### Monte-Carlo Policy Gradient (REINFORCE)

- Update parameters by stochastic gradient ascent
- Using policy gradient theorem
- Using return  $G_t$  as an unbiased sample of  $q_{\pi_{\theta}}(s_t, a_t)$

 $\Delta \theta_t = \alpha \nabla_\theta \log \pi_\theta(s_t, a_t) \, \mathrm{G}_t$ 



[An Introduction to Reinforcement Learning, Sutton and Barto]

# Reducing Variance Using a Critic

-Monte-Carlo policy gradient still has high variance

-We use a critic to estimate the action-value function

 $q_w(s,a) \approx q_{\pi_\theta}(s,a)$ 

- -Actor-critic algorithms maintain two sets of parameters
  - Critic Updates action-value function parameters w
  - Actor Updates policy parameters  $\theta$ , in direction suggested by critic

-Actor-critic algorithms follow an approximate policy gradient

$$\nabla_{\theta} J(\theta) \approx \mathbb{E}_{\pi_{\theta}} [\nabla_{\theta} \log \pi_{\theta}(s, a) q_{w}(s, a)]$$
$$\Delta \theta = \alpha \nabla_{\theta} \log \pi_{\theta}(s, a) q_{w}(s, a)$$

### Estimating the Action-Value Function

- -The critic is solving a familiar problem: policy evaluation (prediction) -How good is policy  $\pi_{\theta}$  for current parameters  $\theta$ ?
- -Could also use e.g., least-squares policy evaluation

# Action-Value Actor-Critic (1/2)

-Simple actor-critic algorithm based on action-value critic

- -Using linear value function approximation  $q_w(s,a) = \varphi(s,a)^T w$ 
  - Critic Updates w by linear TD(0)
  - Actor Updates  $\theta$  by policy gradient

# Action-Value Actor-Critic (2/2)

One-step Actor–Critic (episodic), for estimating  $\pi_{\theta} \approx \pi_*$ 

Input: a differentiable policy parameterization  $\pi(a|s, \theta)$ Input: a differentiable state-value function parameterization  $\hat{v}(s, \mathbf{w})$ Parameters: step sizes  $\alpha^{\theta} > 0, \ \alpha^{\mathbf{w}} > 0$ Initialize policy parameter  $\boldsymbol{\theta} \in \mathbb{R}^{d'}$  and state-value weights  $\mathbf{w} \in \mathbb{R}^{d}$  (e.g., to **0**) Loop forever (for each episode): Initialize S (first state of episode)  $I \leftarrow 1$ Loop while S is not terminal (for each time step):  $A \sim \pi(\cdot | S, \theta)$ Take action A, observe S', R(if S' is terminal, then  $\hat{v}(S', \mathbf{w}) \doteq 0$ )  $\delta \leftarrow R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})$  $\mathbf{w} \leftarrow \mathbf{w} + \alpha^{\mathbf{w}} \delta \nabla \hat{v}(S, \mathbf{w})$  $\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} + \alpha^{\boldsymbol{\theta}} I \delta \nabla \ln \pi (A|S, \boldsymbol{\theta})$  $I \leftarrow \gamma I$  $S \leftarrow S'$ 

[An Introduction to Reinforcement Learning, Sutton and Barto]

### Approximation with Deep Networks

- -So far, the feature representation was typically "fixed"
- –The parametrised functions  $\hat{v}(s; w) / \pi_{\theta}(s, a)$  were linear mappings of input features
- More complicated non-linear mappings are needed to generalize to more complex domains
- -Popular choice is to use deep neural networks
  - Known to discover useful feature representation tailored to the specific task
  - We can leverage extensive research on architectures and optimisation from Supervised Learning

# Batch Reinforcement Learning

- -Gradient descent is simple and appealing
- -But it is not sample efficient
- -Batch methods seek to find the best fitting value function
- -Given the agent's experience ("training data")

### Stochastic Gradient Descent with Experience Replay

-Given experience consisting of (*state*, *value*) pairs

$$\mathcal{D} = \left\{ \langle s_1, \vee_1^{\pi} \rangle, \left\langle \langle s_2, \vee_2^{\pi} \rangle \right\rangle, \dots, \left\langle \langle s_T, \vee_T^{\pi} \rangle \right\rangle \right\}$$

-Repeat:

1. Sample state, value from experience

 $\langle s, \vee^{\pi} \rangle \sim \mathcal{D}$ 

2. Apply stochastic gradient descent update

$$\Delta w = \alpha(\vee_{\pi}(s) - \hat{\vee}(s; w))\nabla_{w}\hat{\vee}(s; w)$$
  
Converges to least squares solution  
$$w^{\pi} = \underset{w}{\operatorname{argmin}} LS(w)$$

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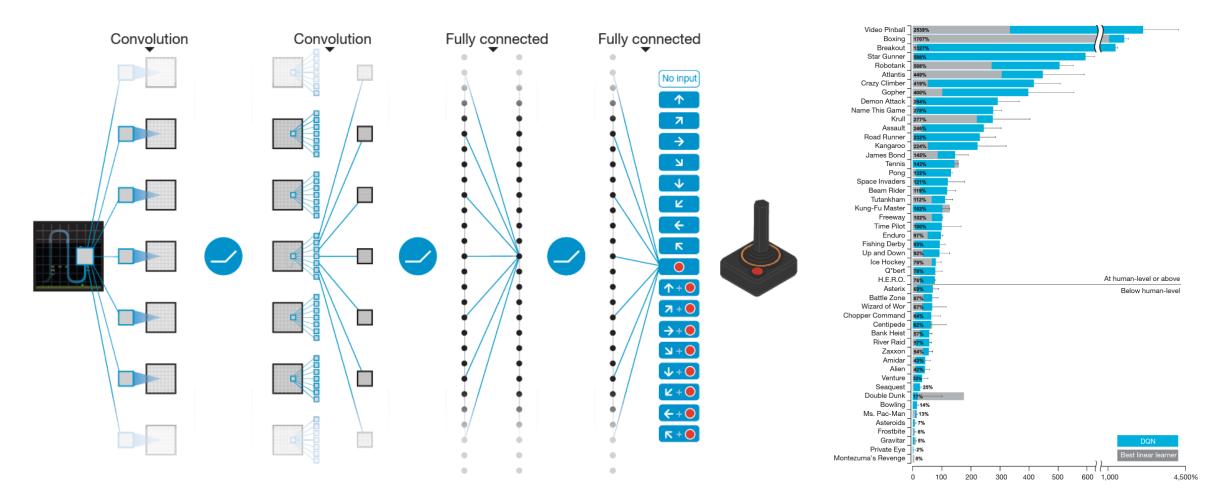
### Experience Replay in Deep Q-Networks (DQN)

-Example of DQN that uses experience replay and fixed Q-targets

- Take action  $a_t$  according to  $\varepsilon$ -greedy policy
- Store transition  $(s_t, a_t, r_{t+1}, s_{t+1})$  in replay memory  $\mathcal{D}$
- Sample random mini-batch of transitions (s, a, r, s') from  $\mathcal{D}$
- Compute Q-learning targets w.r.t. the parameters of a DQN  $w^-$
- Optimize MSE between Q-network and Q-learning targets

$$\mathcal{L}_{i}(w_{i}) = \mathbb{E}_{s,a,r,s'\sim\mathcal{D}_{i}}\left[\left(r + \gamma \max_{a'} Q(s',a';w_{i}^{-}) - Q(s,a;w_{i})\right)^{2}\right]$$

### DQN in Atari



### **Continuous Action Spaces**

-Vanilla DQN can't be used for continuous action spaces (CAS)

- Too many outputs to parametrize with a neural network
- Discretization would be needed (suboptimal)
- -Deep Deterministic Policy Gradient (DDPG) extends DQN to CAS
  - DDPG is an Actor Critic algorithm with critic  $\pi_{\theta}$  and actor  $Q_w$  networks
  - Policy is deterministic, noise added for exploration  $a_t = \pi_{\theta}(s_t) + \varepsilon$
  - Update the critic with the actor as max and then the actor to maximize the critic

$$\mathcal{L}_i(w_i) = \mathbb{E}_{s,a,r,s' \sim \mathcal{D}_i}[(r + \gamma Q(s', \pi(s'; \theta_i^-); w_i^-) - Q(s, a; w_i))^2]$$

$$\mathcal{L}_i(\theta_i) = \mathbb{E}_{s \sim \mathcal{D}_i}[Q(s, \pi(s; \theta); w_i)]$$

### DDPG

#### Algorithm 1 Deep Deterministic Policy Gradient

- 1: Input: initial policy parameters  $\theta$ , Q-function parameters  $\phi$ , empty replay buffer  $\mathcal{D}$
- 2: Set target parameters equal to main parameters  $\theta_{targ} \leftarrow \theta$ ,  $\phi_{targ} \leftarrow \phi$

#### 3: repeat

- 4: Observe state s and select action  $a = \operatorname{clip}(\mu_{\theta}(s) + \epsilon, a_{Low}, a_{High})$ , where  $\epsilon \sim \mathcal{N}$
- 5: Execute a in the environment
- 6: Observe next state s', reward r, and done signal d to indicate whether s' is terminal
- 7: Store (s, a, r, s', d) in replay buffer  $\mathcal{D}$
- 8: If s' is terminal, reset environment state.
- 9: if it's time to update then
- 10: for however many updates do
- 11: Randomly sample a batch of transitions,  $B = \{(s, a, r, s', d)\}$  from  $\mathcal{D}$
- 12: Compute targets

$$y(r, s', d) = r + \gamma(1 - d)Q_{\phi_{\text{targ}}}(s', \mu_{\theta_{\text{targ}}}(s'))$$

13: Update Q-function by one step of gradient descent using

$$\nabla_{\phi} \frac{1}{|B|} \sum_{(s,a,r,s',d) \in B} (Q_{\phi}(s,a) - y(r,s',d))^2$$

14: Update policy by one step of gradient ascent using

$$\nabla_{\theta} \frac{1}{|B|} \sum_{s \in B} Q_{\phi}(s, \mu_{\theta}(s))$$

15: Update target networks with

$$\phi_{\text{targ}} \leftarrow \rho \phi_{\text{targ}} + (1 - \rho) \phi$$
$$\theta_{\text{targ}} \leftarrow \rho \theta_{\text{targ}} + (1 - \rho) \theta$$

16: end for
17: end if
18: until convergence