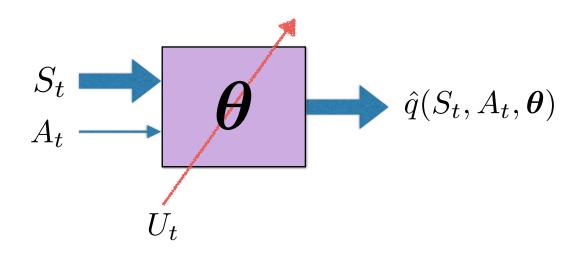
Sequential decision making Control: Deep Q-Learning (DQN) and Eligibility Trace

Value function approximation (VFA) for control



Recall: Different Targets

- Monte Carlo: $G_t \doteq R_{t+1} + \gamma R_{t+2} + \gamma^2 R_{t+3} + \dots + \gamma^{T-t-1} R_T$
- TD: $G_t^{(1)} \doteq R_{t+1} + \gamma V_t(S_{t+1})$
 - Use V_t to estimate remaining return
- *n*-step TD:
 - **2 step return:** $G_t^{(2)} \doteq R_{t+1} + \gamma R_{t+2} + \gamma^2 V_t(S_{t+2})$
 - *n*-step return: $G_t^{(n)} \doteq R_{t+1} + \gamma R_{t+2} + \gamma^2 + \dots + \gamma^{n-1} R_{t+n} + \gamma^n V_t(S_{t+n})$ $G_t^{(n)} \doteq G_t \text{ if } t + n \ge T$

Recall: Stochastic Gradient Descent (SGD)

```
General SGD: \boldsymbol{\theta} \leftarrow \boldsymbol{\theta} - \alpha \nabla_{\boldsymbol{\theta}} \operatorname{Error}_{t}^{2}

For VFA: \leftarrow \boldsymbol{\theta} - \alpha \nabla_{\boldsymbol{\theta}} \left[ \operatorname{Target}_{t} - \hat{v}(S_{t}, \boldsymbol{\theta}) \right]^{2}

Chain rule: \leftarrow \boldsymbol{\theta} - 2\alpha \left[ \operatorname{Target}_{t} - \hat{v}(S_{t}, \boldsymbol{\theta}) \right] \nabla_{\boldsymbol{\theta}} \left[ \operatorname{Target}_{t} - \hat{v}(S_{t}, \boldsymbol{\theta}) \right]

Semi-gradient: \leftarrow \boldsymbol{\theta} + \alpha \left[ \operatorname{Target}_{t} - \hat{v}(S_{t}, \boldsymbol{\theta}) \right] \nabla_{\boldsymbol{\theta}} \hat{v}(S_{t}, \boldsymbol{\theta})
```

Different RL algorithms provide different targets!

But share the "semi-gradient" aspect

(Semi-)gradient methods carry over to control in the usual on-policy GPI way

- Always learn the action-value function of the current policy
- Always act near-greedily wrt the current action-value estimates
- The learning rule is:

$$\boldsymbol{\theta}_{t+1} \doteq \boldsymbol{\theta}_t + \alpha \begin{bmatrix} U_t - \hat{q}(S_t, A_t, \boldsymbol{\theta}_t) \end{bmatrix} \nabla \hat{q}(S_t, A_t, \boldsymbol{\theta}_t)$$
 update target, e.g. $U_t = G_t$ (MC)
$$U_t = R_{t+1} + \gamma \hat{q}(S_{t+1}, A_{t+1}, \boldsymbol{\theta}_t) \text{ (Sarsa)}$$

$$U_t = R_{t+1} + \gamma \sum_a \pi(a|S_{t+1}) \hat{q}(S_{t+1}, a, \boldsymbol{\theta}_t) \qquad U_t = \sum_{s', r} p(s', r|S_t, A_t) \Big[r + \gamma \sum_{a'} \pi(a'|s') \hat{q}(s', a', \boldsymbol{\theta}_t) \Big] \text{ (DP)}$$
 (Expected Sarsa)

(Semi-)gradient methods carry over to control

$$\boldsymbol{\theta}_{t+1} \doteq \boldsymbol{\theta}_t + \alpha \Big[U_t - \hat{q}(S_t, A_t, \boldsymbol{\theta}_t) \Big] \nabla \hat{q}(S_t, A_t, \boldsymbol{\theta}_t)$$

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

Input: a differentiable function $\hat{q}: \mathcal{S} \times \mathcal{A} \times \mathbb{R}^n \to \mathbb{R}$ Initialize value-function weights $\boldsymbol{\theta} \in \mathbb{R}^n$ arbitrarily (e.g., $\boldsymbol{\theta} = \mathbf{0}$)

Repeat (for each episode):

 $S, A \leftarrow \text{initial state}$ and action of episode (e.g., ε -greedy)

Repeat (for each step of episode):

Take action A, observe R, S'

If S' is terminal:

$$\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} + \alpha [R - \hat{q}(S, A, \boldsymbol{\theta})] \nabla \hat{q}(S, A, \boldsymbol{\theta})$$

Go to next episode

Choose A' as a function of $\hat{q}(S', \cdot, \boldsymbol{\theta})$ (e.g., ε -greedy)

$$\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} + \alpha \left[R + \gamma \hat{q}(S', A', \boldsymbol{\theta}) - \hat{q}(S, A, \boldsymbol{\theta}) \right] \nabla \hat{q}(S, A, \boldsymbol{\theta})$$

 $S \leftarrow S'$

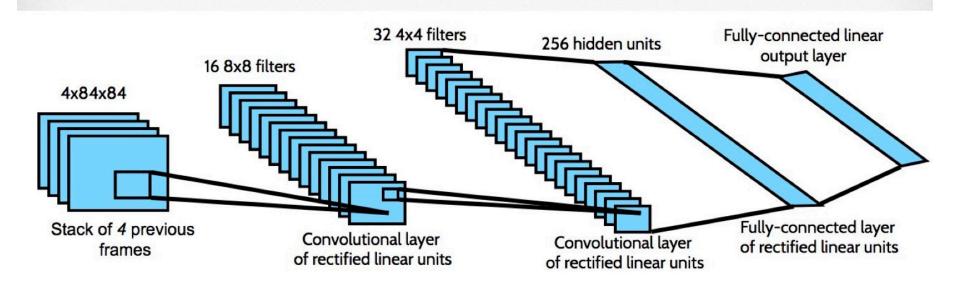
$$A \leftarrow A'$$

DQN

(Mnih, Kavukcuoglu, Silver, et al., Nature 2015)

- Learns to play video games from raw pixels, simply by playing
- Can learn Q function by Q-learning

$$\Delta \boldsymbol{w} = \alpha \left(R_{t+1} + \gamma \max_{a} Q(S_{t+1}, a; \boldsymbol{w}) - Q(S_{t}, A_{t}; \boldsymbol{w}) \right) \nabla_{\boldsymbol{w}} Q(S_{t}, A_{t}; \boldsymbol{w})$$



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- Core components of DQN include:
 - Target networks (Mnih et al. 2015)

$$\Delta \boldsymbol{w} = \alpha \left(R_{t+1} + \gamma \max_{\boldsymbol{a}} Q(S_{t+1}, \boldsymbol{a}; \boldsymbol{w}^{-}) - Q(S_{t}, A_{t}; \boldsymbol{w}) \right) \nabla_{\boldsymbol{w}} Q(S_{t}, A_{t}; \boldsymbol{w})$$

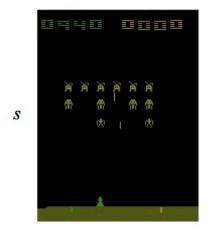
Experience replay (Lin 1992): replay previous tuples (s, a, r, s')

Target Network Intuition

(Slide credit: Vlad Mnih)

- Changing the value of one action will change the value of other actions and similar states.
- The network can end up chasing its own tail because of bootstrapping.
- Somewhat surprising fact bigger networks are less prone to this because they alias less.

$$L_i(\theta_i) = \mathbb{E}_{s,a,s',r \sim D} \left(\underbrace{r + \gamma \, \max_{a'} Q(s', a'; \boldsymbol{\theta}_i^-)}_{\text{target}} - Q(s, a; \theta_i) \right)^2$$





DQN

(Mnih, Kavukcuoglu, Silver, et al., Nature 2015)

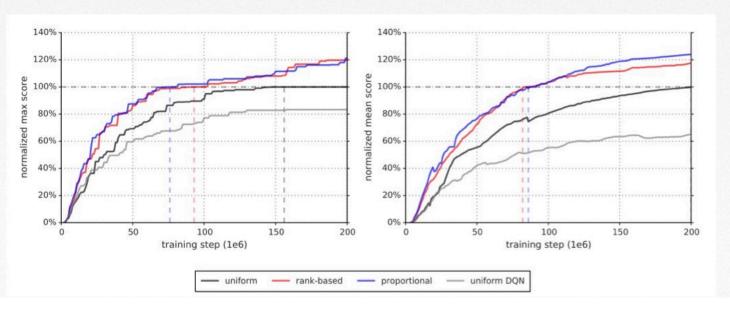
- Many later improvements to DQN
 - Double Q-learning (van Hasselt 2010, van Hasselt et al. 2015)
 - Prioritized replay (Schaul et al. 2016)
 - Dueling networks (Wang et al. 2016)
 - Asynchronous learning (Mnih et al. 2016)
 - Adaptive normalization of values (van Hasselt et al. 2016)
 - Better exploration (Bellemare et al. 2016, Ostrovski et al., 2017, Fortunato, Azar, Piot et al. 2017)
 - Distributional losses (Bellemare et al. 2017)
 - o Multi-step returns (Mnih et al. 2016, Hessel et al. 2017)
 - o ... many more ...

Prioritized Experience Replay

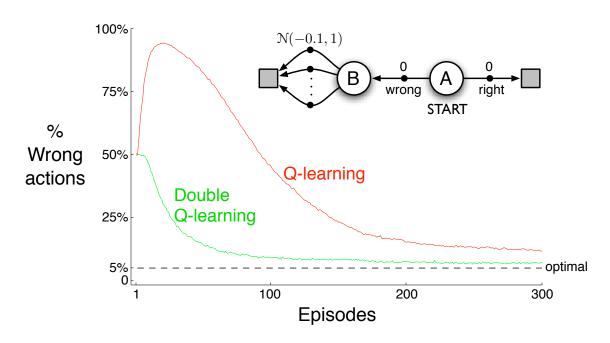
"Prioritized Experience Replay", Schaul et al. (2016)

Idea: Replay transitions in proportion to TD error:

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



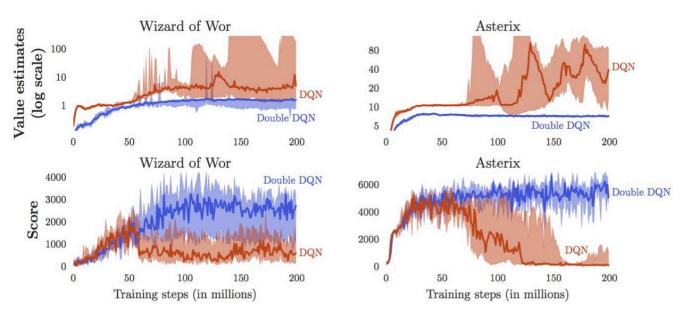
Recall: Double DQN



Double Q-learning:

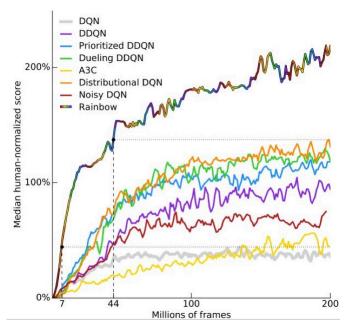
$$Q_1(S_t, A_t) \leftarrow Q_1(S_t, A_t) + \alpha \left[R_{t+1} + \gamma Q_2(S_{t+1}, \arg\max_{a} Q_1(S_{t+1}, a)) - Q_1(S_t, A_t) \right]$$

Double DQN



cf. van Hasselt et al, 2015)

Which DQN improvements

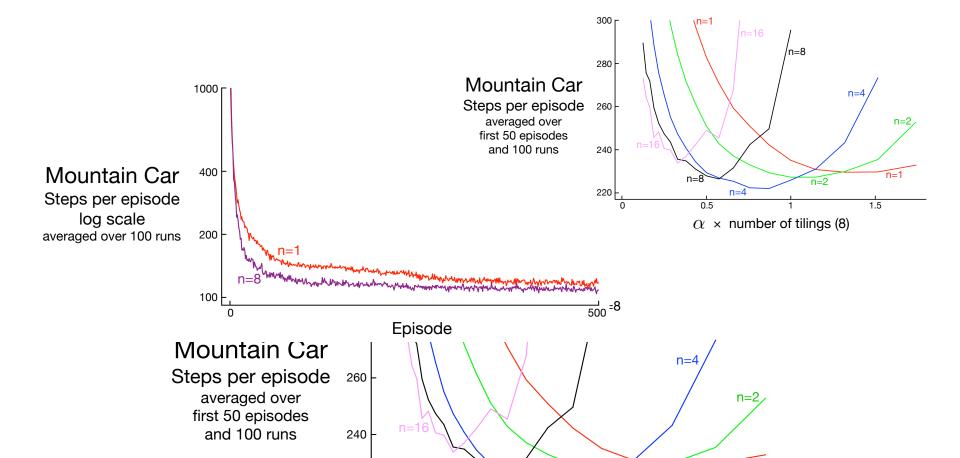


Rainbow model, Hessel et al, 2017)

Deep n-step SARSA

```
Episodic semi-gradient n-step Sarsa for estimating \hat{q} \approx q_* or q_{\pi}
Input: a differentiable action-value function parameterization \hat{q}: \mathbb{S} \times \mathcal{A} \times \mathbb{R}^d \to \mathbb{R}
Input: a policy \pi (if estimating q_{\pi})
Algorithm parameters: step size \alpha > 0, small \varepsilon > 0, a positive integer n
Initialize 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, \text{ and } R_t) can take their index mod n+1
Loop for each episode:
    Initialize and store S_0 \neq \text{terminal}
    Select and store an action A_0 \sim \pi(\cdot|S_0) or \varepsilon-greedy wrt \hat{q}(S_0,\cdot,\mathbf{w})
    T \leftarrow \infty
    Loop for t = 0, 1, 2, ...:
        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 \varepsilon-greedy wrt \hat{q}(S_{t+1}, \cdot, \mathbf{w})
        \tau \leftarrow t - n + 1 (\tau is the time whose estimate is being updated)
        If \tau > 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})
             \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
```





Eligibility traces are

- Another way of interpolating between MC and TD methods
- A way of implementing *compound* λ -return targets
- A basic mechanistic idea a short-term, fading memory
- A new style of algorithm development/ analysis

Recall *n*-step targets

- For example, in the episodic case, with linear function approximation:
 - 2-step target: $G_t^{(2)} \doteq R_{t+1} + \gamma R_{t+2} + \gamma^2 \theta_{t+1}^{\top} \phi_{t+2}$
 - *n*-step target: $G_t^{(n)} \doteq R_{t+1} + \dots + \gamma^{n-1} R_{t+n} + \gamma^n \boldsymbol{\theta}_{t+n-1}^{\top} \boldsymbol{\phi}_{t+n}$ with $G_t^{(n)} \doteq G_t$ if $t+n \geq T$

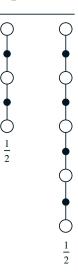
Any set of update targets can be

• For example, half a 2-step plus half a 4step 1 = (2) 1 = (4)

$$U_t = \frac{1}{2}G_t^{(2)} + \frac{1}{2}G_t^{(4)}$$

- Called a compound backup
 - Draw each component
 - Label with the weights for that

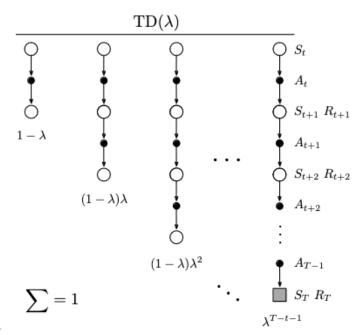
A compound backup



The λ -return is a compound update target

- The λ -return a target that averages all n-step targets
 - each weighted by λ^{n-1}

$$G_t^{\lambda} \doteq (1 - \lambda) \sum_{n=1}^{\infty} \lambda^{n-1} G_{t:t+n}.$$



 $G_{t:t+n} \doteq R_{t+1} + \gamma R_{t+2} + \dots + \gamma^{n-1} R_{t+n} + \gamma^n \hat{v}(S_{t+n}, \mathbf{w}_{t+n-1}), \quad 0 \le t \le T - n$

Relation to TD(0) and MC

• The λ -return can be rewritten as:

$$G_t^{\lambda} = (1-\lambda)\sum_{n=1}^{T-t-1}\lambda^{n-1}G_t^{(n)} + \lambda^{T-t-1}G_t$$
 Until termination After termination

• If $\lambda = 1$, you get the MC target:

$$G_t^{\lambda} = (1-1)\sum_{n=1}^{T-t-1} 1^{n-1} G_t^{(n)} + 1^{T-t-1} G_t = G_t$$

• If $\lambda = 0$, you get the TD(0) target:

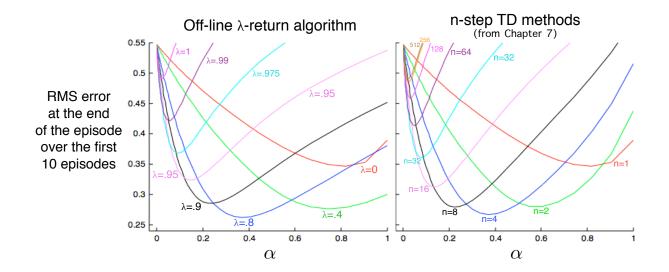
$$G_t^{\lambda} = (1-0) \sum_{n=1}^{T-t-1} 0^{n-1} G_t^{(n)} + 0^{T-t-1} G_t = G_t^{(1)}$$
 21

The off-line λ -return "algorithm"

Wait until the end of the episode (offline)

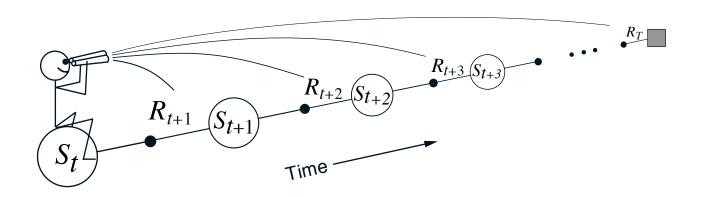
$$\boldsymbol{\theta}_{t+1} \doteq \boldsymbol{\theta}_t + \alpha \left[\boldsymbol{G}_t^{\lambda} - \hat{q}(S_t, A_t, \boldsymbol{\theta}_t) \right] \nabla \hat{q}(S_t, A_t, \boldsymbol{\theta}_t), \quad t = 0, \dots, T-1$$

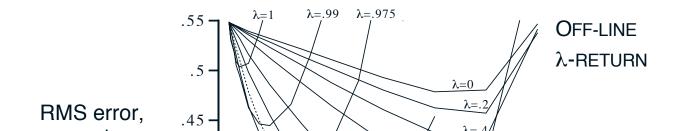
The λ -return alg performs similarly to



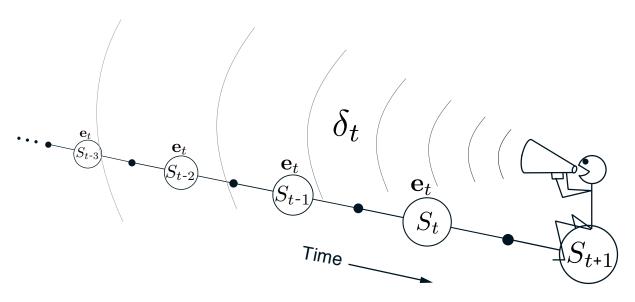
Intermediate λ is best (just like intermediate n is best) λ -return slightly better than n-step

The forward view looks forward from the state being updated to future states and rewards





The backward view looks back to the recently visited states (marked by eligibility traces)



- Shout the TD error backwards
- The traces fade with temporal distance by $\gamma\lambda$

Eligibility traces (mechanism)

- The forward view was for theory
- The backward view is for mechanism same shape as θ

$$\mathbf{e}_t \in \mathbb{R}^{n'} \geq \mathbf{0}$$

- New memory vector called *eligibility trace*
 - On each step, decay each component by $\gamma\lambda$ and increment the trace for the current state by 1
 - Accumulating trace

$$\mathbf{e}_0 \doteq \mathbf{0},$$

 $\mathbf{e}_t \doteq \nabla \hat{v}(S_t, \boldsymbol{\theta}_t) + \gamma \lambda \mathbf{e}_{t-1}$

The Semi-gradient $TD(\lambda)$ algorithm

$$\theta_{t+1} \doteq \theta_t + \alpha \delta_t \mathbf{e}_t$$

$$\delta_t \doteq R_{t+1} + \gamma \hat{v}(S_{t+1}, \theta_t) - \hat{v}(S_t, \theta_t)$$

$$\mathbf{e}_0 \doteq \mathbf{0},$$

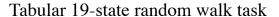
$$\mathbf{e}_t \doteq \nabla \hat{v}(S_t, \theta_t) + \gamma \lambda \mathbf{e}_{t-1}$$

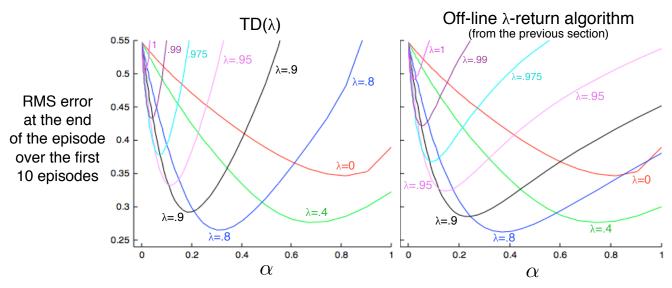
Online $TD(\lambda)$

Semi-gradient TD(λ) 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} such that \hat{v}(\text{terminal}, \cdot) = 0
Algorithm parameters: step size \alpha > 0, trace decay rate \lambda \in [0, 1]
Initialize value-function weights \mathbf{w} arbitrarily (e.g., \mathbf{w} = \mathbf{0})
Loop for each episode:
    Initialize S
                                                                                           (a d-dimensional vector)
    z \leftarrow 0
    Loop for each step of episode:
         Choose A \sim \pi(\cdot|S)
         Take action A, observe R, S'
         \mathbf{z} \leftarrow \gamma \lambda \mathbf{z} + \nabla \hat{v}(S, \mathbf{w})
        \delta \leftarrow R + \gamma \hat{v}(S', \mathbf{w}) - \hat{v}(S, \mathbf{w})
         \mathbf{w} \leftarrow \mathbf{w} + \alpha \delta \mathbf{z}
         S \leftarrow S'
    until S' is terminal
```

TD(λ) performs similarly to offline λ -





Can we do better? Can we update online?

Conclusions

- Value-function approximation by stochastic gradient descent enables RL to be applied to arbitrarily large state spaces
- Most algorithms just carry over the targets from the tabular case
- With bootstrapping (TD), we don't get true gradient descent methods
 - this complicates the analysis
 - but the linear, on-policy case is still guaranteed convergent
 - and learning is still much faster