Reinforcement Learning

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Deep Neural Networks

Use of Neural Networks for regression



Two possible approaches for function approximation:

- Incremental:
 - Pro: Learning on-line
 - Cons: No convergence due to (a) Data not i.i.d., that can lead to catastrophic forgetting, and (b) Moving target problem

Two possible approaches for function approximation:

- Incremental:
 - Pro: Learning on-line
 - Cons: No convergence due to (a) Data not i.i.d., that can lead to catastrophic forgetting, and (b) Moving target problem
- Batch Learning:
 - Cons: Learn from collected dataset (not own experience)
 - Pro: Better convergence

Fitted Q-learning

Given \mathcal{D} of size T with examples $(s_t, a_t, r_{t+1}, s_{t+1})$, and regression algorithm, set N to zero and $Q_N(s, a) = 0$ for all a and s repeat

 $\begin{array}{l} N \leftarrow N+1 \\ \text{Build training set } TS = \{\langle (s_t, a_t), r_{t+1} + \gamma \max_a Q_N(s_{t+1}, a) \rangle \}_{t=1}^T \\ Q_{N+1} \leftarrow \text{ regression algorithm on TS} \\ \textbf{until } Q_N \approx Q_{N+1} \text{ or } N > \text{limit} \\ \textbf{return } \pi \text{ based on greedy evaluation of } Q_N \end{array}$

Neural Fitted Q-learning: Wrong version. Why?

Initialize weights θ for NN for regression Collect \mathcal{D} of size T with examples $(s_t, a_t, r_{t+1}, s_{t+1})$ **repeat** Sample \mathcal{B} mini-batch of \mathcal{D} $\theta \leftarrow \theta - \alpha \sum_{t \in \mathcal{B}} \frac{\partial Q_{\theta}}{\partial \theta}(s_t, a_t) (Q_{\theta}(s_t, a_t) - [r_{t+1} + \gamma \max_{a'} Q_{\theta}(s_{t+1}, a')])$ **until** convergence on learning or maximum number of steps **return** π based on greedy evaluation of Q_{θ}

- Does not work well
- It's not a Batch method. Can you see why?

Neural Fitted Q-learning

```
Initialize weights \theta for NN for regression

Collect \mathcal{D} of size T with examples (s_t, a_t, r_{t+1}, s_{t+1})

repeat

\theta' \leftarrow \theta

repeat

Sample \mathcal{B} mini-batch of \mathcal{D}

\theta \leftarrow \theta - \alpha \sum_{t \in \mathcal{B}} \frac{\partial Q_{\theta}}{\partial \theta}(s_t, a_t) (Q_{\theta}(s_t, a_t) - [r_{t+1} + \gamma \max_{a'} Q_{\theta'}(s_{t+1}, a')])

until convergence on learning or maximum number of steps

until maximum limit iterations

return \pi based on greedy evaluation of Q'_{\theta}
```

• Notice target does not change during supervised regression

Neural Fitted Q-learning: Another version

- That works, however the update of parameters is not smooth
- Alternative version to avoid moving target

Fitted Q-learning avoiding moving target

Initialize weights θ for NN for regression Collect \mathcal{D} of size T with examples $(s_t, a_t, r_{t+1}, s_{t+1})$ **repeat** Sample \mathcal{B} mini-batch of \mathcal{D} $\theta \leftarrow \theta - \alpha \sum_{t \in \mathcal{B}} \frac{\partial Q_{\theta}}{\partial \theta}(s_t, a_t) - (Q_{\theta}(s_t, a_t) - [r_{t+1} + \gamma \max_{a'} Q_{\theta'}(s_{t+1}, a')])$ $\theta' \leftarrow \tau \theta' + (1 - \tau)\theta$ **until** maximum limit iterations **return** π based on greedy evaluation of Q'_{θ}

• Value of τ close to one (f.i. $\tau = 0.999$) reduces the "speed" of the moving target.

- So now, we have learning stabilized just any batch method but using NN.
- \bullet However, now there is the problem of dependence of dataset $\mathcal{D}.$ How we obtain the data?
- Data can be obtained using a random policy, but we want to minimize error on states visited by the policy!

$$L(heta) = \mathbb{E}_{\pi}\left[(V^{\pi}(s) - V_{ heta}(s))^2
ight] = \sum_{s \in \mathcal{S}} \mu^{\pi}(s) \left[V^{\pi}(s) - V_{ heta}(s)
ight]^2$$

where $\mu^{\pi}(s)$ is the time spent in state s while following π

- Data should be generated by the policy
- But it also has to be probabilistic (to ensure exploration)
- \bullet So, collect data using the policy and add them to ${\cal D}$
- Also remove old data from \mathcal{D} .
 - Limit the size of the set
 - Remove examples obtained using old policies
- So, collect data using a *buffer* of limited size (we call replay buffer).

Batch Q-learning with replay buffer and target network

Initialize weights θ for NN for regression

Collect D of size T with examples $(s_t, a_t, r_{t+1}, s_{t+1})$ using random policy

repeat

 $\theta' \leftarrow \theta$

repeat

Collect *M* experiences following ϵ -greedy procedure and add them to **buffer** \mathcal{D}

repeat

Sample ${\mathcal B}$ mini-batch of ${\mathcal D}$

 $\theta \leftarrow \theta - \alpha \sum_{t \in \mathcal{B}} \frac{\partial Q_{\theta}}{\partial \theta}(s_t, a_t) \left(Q_{\theta}(s_t, a_t) - [r_{t+1} + \gamma \max_{a'} Q_{\theta'}(s_{t+1}, a')] \right)$

until maximum number of steps K

until maximum number of iterations N

until maximum limit iterations

return π based on greedy evaluation of Q'_{θ}

DQN algorithm (Mnih, et al. 2015)

• Deep Q-Network algorithm breakthrough

- ► In 2015, Nature published DQN algorithm.
- It takes profit of "then-recent" Deep Neural Networks and, in particular, of Convolutional NNs so successful for vision problems
- Applied to Atari games directly from pixels of the screen (no hand made representation of the problem)
- Very successful on a difficult task, surpassing in some cases human performance
- It is basically the previous algorithm with K = 1, and M = 1 that is applied on the current state.
- It goes back to incremental learning

DQN algorithm

Initialize weights θ for NN for regression Set *s* to initial state, and *k* to zero

repeat

Choose *a* from *s* using policy π_{θ} derived from Q_{θ} (e.g., ϵ -greedy) $k \leftarrow k + 1$ Execute action *a*, observe *r*, *s'*, and add $\langle s, a, r, s' \rangle$ to buffer \mathcal{D} Sample \mathcal{B} mini-batch of \mathcal{D} $\theta \leftarrow \theta - \alpha \sum_{t \in \mathcal{B}} \frac{\partial Q_{\theta}}{\partial \theta}(s_t, a_t) (Q_{\theta}(s_t, a_t) - [r_{t+1} + \gamma \max_{a'} Q_{\theta'}(s_{t+1}, a')])$ if k==N then $\theta' \leftarrow \theta$ $k \leftarrow 0$ end if until maximum limit iterations return π based on greedy evaluation of Q'_{θ}

• End-to-end learning of values Q(s; a) from pixels:

State: Input state s is stack of raw pixels from last 4 frames **Actions:** Output is Q(s, a) value for each of 18 joystick/button positions

Reward: Reward is direct change in score for that step

- Network architecture and hyper-parameters **fixed across all games**, No tuning!
- Clipping reward -1,0,1 to avoid problem of different magnitudes of score in each game







• What is the effect of each trick on Atari games?

DQN

	Q-learning	Q-learning	Q-learning	Q-learning
			+ Replay	+ Replay
		+ Target Q		+ Target Q
Breakout	3	10	241	317
Enduro	29	142	831	1006
River Raid	1453	2868	4103	7447
Seaquest	276	1003	823	2894
Space Invaders	302	373	826	1089

Overestimates: Double Q-learning

Double Q-learning (Hasselt, et al. 2015)

• Problem of overestimation of Q values.

• We use "max" operator to compute the target in the minimization of:

$$L(s,a) = (Q(s,a) - (r + \gamma \max_{a'} Q(s',a')))^2$$

- Surprisingly here is a problem.
 - **1** Suppose Q(s', a') is 0 for all actions, so Q(s, a) should be r.
 - But γ max_a, Q(s', a') ≥ 0 because random initialization and use of the max operator.
 - **③** So estimation $Q(s, a) \ge r$, overestimating true value
 - All this because for max operator:

$$\mathbb{E}[\max_{a'} Q(s',a')] \geq \max_{a'} \mathbb{E}[Q(s',a')]$$

• This overestimation is propagated to other states.

Double Q-learning

- Solution (Hasselt, 2010): Train 2 action-value functions: Q_A and Q_B , and compute argmax with the other network
- Do Q-learning on both, but
 - never on the same time steps (Q_A and Q_B are independent)
 - pick Q_A or Q_B at random to be updated on each step
- Notice that:

$$r + \gamma \max_{a'} Q(s, a') = r + \gamma Q(s, \arg \max_{a'} Q(s', a'))$$

• When updating one network, use the values of the other network:

$$egin{aligned} Q_A(s,a) \leftarrow r + \gamma Q_B(s,rgmax Q_A(s',a')) \ Q_B(s,a) \leftarrow r + \gamma Q_A(s,rgmax Q_B(s',a')) \ a' \end{aligned}$$

• Idea is that they should compensate mistakes of each other because they will be independent. When one network overestimate, probably, the other no, so they mutually cancel overestimation

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Reinforcement Learning

Double DQN (Hasselt, et al. 2015)

- In DQN, in fact, we have 2 value functions: $Q_{ heta}$ and $Q_{ heta'}$
- so, no need to add another one:
 - Current Q-network θ is used to select actions
 - Older Q-network θ' is used to evaluate actions
- Update in Double-DQN (Hasselt, et al. 2015):

$$Q_{\theta}(s, a) \leftarrow r + \gamma \underbrace{\overbrace{Q_{\theta'}(s, \arg\max_{a'} Q_{\theta}(s', a'))}^{\text{Action Evaluation}}}_{\text{Action Selection}}$$

• Works well in practice.

Prioritized Experience Replay

- Idea: sample transitions from replay buffer more cleverly
- Those states with poorer estimation in buffer will be selected with preference for update
- We will set probability for every transition. Let's use the absolute value of TD-error of transition as a probability!

$$p_i = |\mathsf{TD}\operatorname{-error}_i| = |Q_{ heta'}(s_i, a_i) - (r_i + \gamma Q_{ heta'}(s_i, rg\max_{a'} Q_{ heta}(s_{i+1}, a'))|$$

$$P(i) = \frac{p_i^{\alpha}}{\sum_k p_k^{\alpha}}$$

where P(i) is probability of selecting sample *i* for the mini-batch, and $\alpha \ge 0$ is a new parameter ($\alpha = 0$ implies uniform probability)

Prioritized Experience Replay (Schaul, et al. 2016)

• Do you see any problem?

- Do you see any problem?
- Now transitions are no i.i.d. and therefore we introduce a bias.
- Solution: we can correct the bias by using **importance-sampling** weights

$$w_i = \left(\frac{1}{N} \cdot \frac{1}{P(i)}\right)^{\beta}$$

- For numerical reasons, we also normalize weights by $\max_i w_i$
- When we put transition into experience replay, we set it to maximal priority p_t = max_{i<t} p_i

Algorithm 1 Double DQN with proportional prioritization

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

Dueling Network Architectures

- Until now, use of generic NN for regression of Q-value function
- Now, specific Deep Architecture specific for RL
- Advantage function definition:

$$A(s,a) = Q(s,a) - V(s)$$

So,

$$Q(s,a) = A(s,a) + V(s)$$

 Intuitively, Advantage function is relative measure of importance of each action

• Dueling network:



• Intuitive idea is that now we don't learn Q(s, a) independently but share part that is V(s) that improves generalization across actions

- We have now 3 sets of parameters:
 - θ : Usual weights of NN until red section
 - β : Weights to compute V(s)
 - α : Weights to compute A(s, a)
- Green part computes A(s, a) + V(s)



- However, there is a problem: one extra degree of freedom in targets!
- Example:



• Which is the correct one? Notice that:

$$\pi^*(s) = rgmax_{a \in \mathcal{A}} Q^*(s, a)$$

$$V^*(s) = \max_{a \in \mathcal{A}} Q^*(s, a)$$

• So,

$$\max_{a \in \mathcal{A}} A(s, a) = \max_{a \in \mathcal{A}} (Q(s, a) - V(s))$$
$$= \max_{a \in \mathcal{A}} Q(s, a) - V(s)$$
$$= 0$$

• Of course, for actions $a \neq a^* A(s, a) \leq 0$

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- **Solution**: require max_a A(s, a) to be equal to zero!
- So the Q-function computes as:

$$egin{aligned} \mathcal{Q}_{ heta,lpha,eta}(s,a) &= V_{ heta,eta}(s) + \left(\mathcal{A}_{ heta,lpha}(s,a) - \max_{a'\in\mathcal{A}}\mathcal{A}_{ heta,lpha}(s,a')
ight) \end{aligned}$$

In practice, the authors propose to implement

$$Q_{ heta,lpha,eta}(s, a) = V_{ heta,eta}(s) + \left(eta_{ heta,lpha}(s, a) - rac{1}{|eta|} \sum_{oldsymbol{a}'\in\mathcal{A}} eta_{ heta,lpha}(s, a')
ight)$$

- This variant increases stability of the optimization because now depends on softer measure (*average* instead of *max*)
- Now Q-values loses original semantics, but it not important. The important thing is a *reference* between actions

Multi-step learning

Multi-step learning

- Idea: instead of using TD(0), use n-steps estimators like we described in lecture 2
- In buffer we should store experiences:

$$\left\langle s_t, a_t, r_t, \sum_{i=0}^n \gamma^{i-1} r_{t+1} + \gamma^n \max_{a'} Q_{\theta'}(s_{t+n}, a') \right\rangle$$

Multi-step learning

- Idea: instead of using TD(0), use n-steps estimators like we described in lecture 2
- In buffer we should store experiences:

$$\left\langle s_t, a_t, r_t, \sum_{i=0}^n \gamma^{i-1} r_{t+1} + \gamma^n \max_{a'} Q_{\theta'}(s_{t+n}, a') \right\rangle$$

- Again, there is a problem!
- Only correct when learning on-policy! (not an issue when n = 1)
- How to fix that?
 - Ignore the problem (often works well)
 - Dynamically choose n to get only on-policy data (Store data until not policy action taken)
 - ► Use importance sampling (Munos et al, 2016)

Rainbow: Combining Improvements in Deep Reinforcement Learning

- Idea: Let's try to investigate how each of the different improvements over DQN help to improve performance on the Atari games
- Over DQN, they added the following modifications:
 - Double Q-learning
 - Prioritized replay
 - Dueling networks
 - Multi-step learning
 - Distributional RL
 - Noisy Nets
- They perform an ablation study where over the complete set of improvement, they disable one an measure the performance

Rainbow (Hessel et al. 2017)



Figure 2: Each plot shows, for several agents, the number of games where they have achieved at least a given fraction of human performance, as a function of time. From left to right we consider the 20%, 50%, 100%, 200% and 500% thresholds. On the first row we compare Rainbow to the baselines. On the second row we compare Rainbow to its ablations.

Rainbow (Hessel et al. 2017)



Asynchronous Q-learning

Asynchronous Q-learning (Mnih et al. 2016)

• Idea: Parallelize learning with several workers



• After some time steps, the worker passes gradients to the global network

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Asynchronous Q-learning (Mnih et al. 2016)

Algorithm 1 Asynchronous one-step Q-learning - pseudocode for each actor-learner thread.

// Assume global shared θ , θ^- , and counter T = 0. Initialize thread step counter $t \leftarrow 0$ Initialize target network weights $\theta^- \leftarrow \theta$ Initialize network gradients $d\theta \leftarrow 0$ Get initial state s

repeat

Take action a with ϵ -greedy policy based on $Q(s, a; \theta)$ Receive new state s' and reward r $y = \begin{cases} r & \text{for terminal } s' \\ r + \gamma \max_{a'} Q(s', a'; \theta^{-}) & \text{for non-terminal } s' \end{cases}$ Accumulate gradients wrt θ : $d\theta \leftarrow d\theta + \frac{\partial (y - Q(s, a; \theta))^2}{2}$ s = s' $T \leftarrow T + 1$ and $t \leftarrow t + 1$ if $T \mod I_{target} == 0$ then Update the target network $\theta^- \leftarrow \theta$ end if if $t \mod I_{AsyncUpdate} == 0$ or s is terminal then Perform asynchronous update of θ using $d\theta$. Clear gradients $d\theta \leftarrow 0$. end if until $T > T_{max}$

Faster Deep RL by optimality tightening

Faster Deep RL by optimality tightening (He, et al. 2016)

• From Bellman equation we will obtain a bound for a given Q-value:

$$\begin{aligned} Q(s_t, a_t) &= \mathbb{E}\left[r_{t+1} + \gamma \max_{a'} Q(s_{t+1}, a')\right] \\ &\geq \mathbb{E}\left[r_{t+1} + \gamma r_{t+2} + \gamma^2 r_{t+3} + \ldots + \gamma^k \max_{a'} Q(s_{t+k+1}, a')\right] \\ &\geq \mathbb{E}\left[\underbrace{r_{t+1} + \gamma r_{t+2} + \gamma^2 r_{t+3} + \ldots + \gamma^k Q(s_{t+k+1}, a_{t+k+1})}_{\text{Lower bound } L^{max}}\right] \end{aligned}$$

Faster Deep RL by optimality tightening (He, et al. 2016)

Also we can do this backwards in time. Notice that we had

$$Q(s_t, a_t) \geq r_{t+1} + \gamma r_{t+2} + \gamma^2 r_{t+3} + \ldots + \gamma^k Q(s_{t+k+1}, a_{t+k+1})$$

Q(s_t, a_t)
 Changing indexes

$$Q(s_{t-k-1}, a_{t-k-1}) \geq r_{t-k} + \gamma r_{t-k+1} + \gamma^2 r_{t-k+2} + \ldots + \gamma^k Q(s_t, a_t)$$

• So,

$$Q(s_{t-k-1}, a_{t-k-1}) - r_{t-k} - \gamma r_{t-k+1} - \gamma^2 r_{t-k+2} - \dots \geq \gamma^k Q(s_t, a_t)$$

$$\gamma^{-k} \left[Q(s_{t-k-1}, a_{t-k-1}) - r_{t-k} - \gamma r_{t-k+1} - \gamma^2 r_{t-k+2} - \dots \right] \geq Q(s_t, a_t)$$

• So, finally we have an upper bound:

$$Q(s_t, a_t) \leq \underbrace{\gamma^{-k}Q(s_{t-k}, a_{t-k}) - \gamma^{-k}r_{t-k} - \gamma^{-(k-1)}r_{t-(k-1)} - \dots - \gamma r_{t-1}}_{r_{t-k}}$$

Upper bound Umin

(

Faster Deep RL by optimality tightening (He, et al. 2016)

• And now we can modify our loss function using these bounds:

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

$$L(\theta) = \mathbb{E}\left[(Q_{\theta}(s, a) - y)^{2} + \lambda (L^{max} - Q_{\theta}(s, a))^{2}_{+} + \lambda (Q_{\theta}(s, a) - U^{min})^{2}_{+} \right]$$

where λ is a penalization parameter like C in SVMs

- Eq, minimize the original Bellman error. but also *penalizes breaking the bounds*
- Accelerates over an order of magnitude with respect original DQN in number of experiences needed for learning

Practical tricks

- DQN is more reliable on some tasks than others. Test your implementation on reliable tasks like Pong and Breakout: if it doesn't achieve good scores, something is wrong.
- Large replay buffers improve robustness of DQN, and memory efficiency is key.
- SGD can be slow .. rely on RMSprop (or any new optimizer)
- Convolutional models are more ecient then MLPs
- DQN uses action repeat set to 4 (because fps too high speeds training time)
- DQN receives 4 frames of the game at a time (grayscale)
- ϵ is anealled from 1 to .1

- Patience. Training takes time (roughly hours to day on GPU training to see improvement)
- Always use Double DQN (3 lines of difference from DQN)
- Learning rate scheduling is benecial. Try high learning rates in initial exploration period.
- *Exploration* is key: Try non-standard exploration schedules.
- Always run at least two different seeds when experimenting

Practical tricks

• Bellman errors can be big. Clip gradients or use Huber loss on Bellman error

$$L_{\delta}(y, f(x)) = egin{cases} rac{(y-f(x))^2}{2}, & ext{when } |y-f(x)| \leq \delta \ \delta |y-f(x)| - rac{\delta^2}{2}, & ext{otherwise} \end{cases}$$



- Very large γ or set it to 1 to avoid myopic reward (very large sequences before reward)
- n-steps return helps but careful