Lecture 2.0

Reinforcement Learning: Q-learning

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Recap: Dynamic Programming

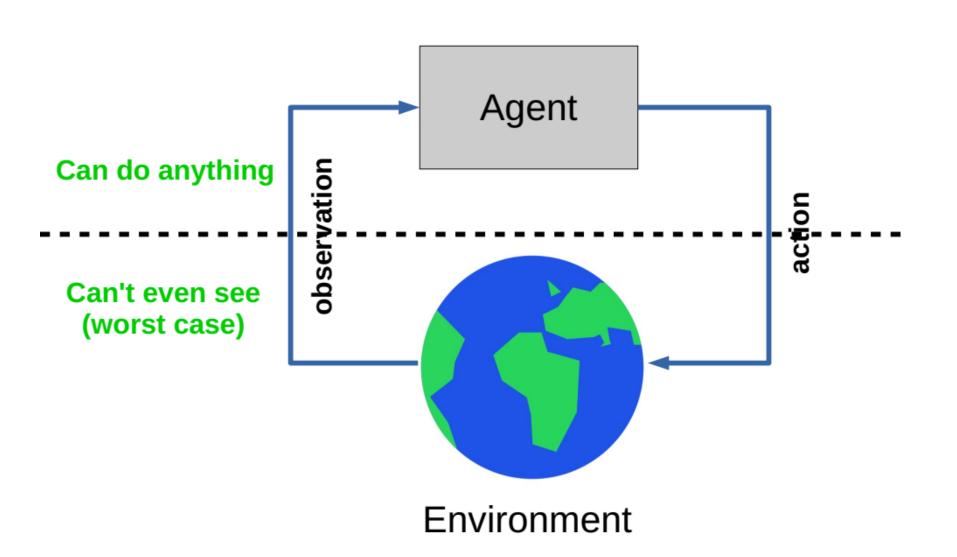
- $v_{\pi}(s), v_{*}(s)$
- If you know v_{*}(s), p(r,s' | s,a) → know optimal policy
- We can learn $v_*(s)$ with Dynamic Programming:

$$v_*(s) = \max_{a} \sum_{r,s'} p(r,s' \mid s,a) [r + \gamma v_*(s')]$$

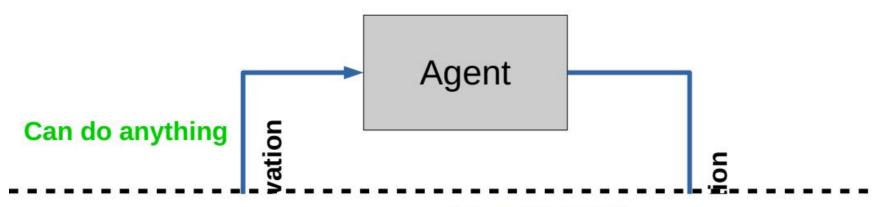
 $q_{\pi}(s, a), q_{*}(s, a)$

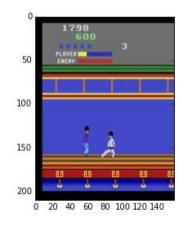
$$q_*(s, a) = \sum_{r,s'} p(r, s' \mid s, a) [r + \gamma \max_{a'} q_*(s', a')]$$

Decision making: reality check



Decision making: reality check











Model-Free Setup

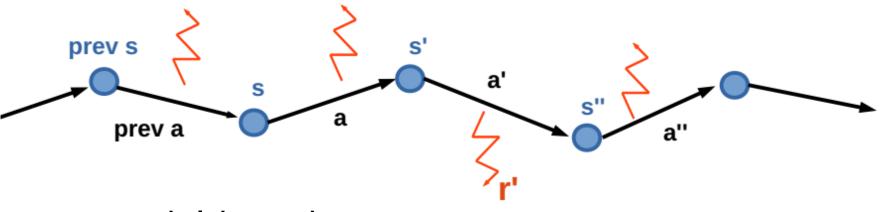
 We don't know internal environment representation, e.g.

$$p(r, s' \mid s, a)$$
 - unknown

What should we do?

Learning from trajectories

 $s_1 -> a_1 -> r_1 -> s_2 -> \dots -> s_n - trajectory$



- Model-based setup:
 - you can apply Dynamic Programming
 - you can plan (!)
- Model-free setup:
 - you can experiment with different actions
 - no guaranties (!!!)

Learning from trajectories

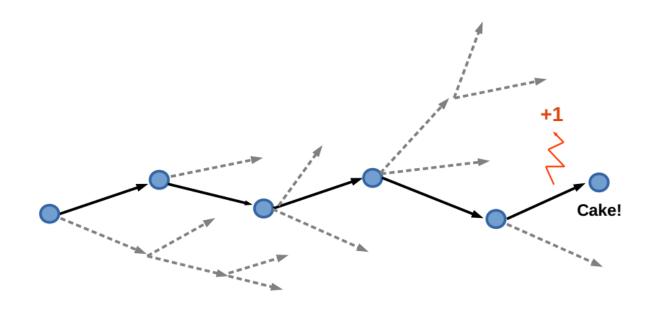
$$s_1 -> a_1 -> r_1 -> s_2 -> \dots -> s_n - trajectory$$

We can sample trajectories (a lot of trajectories!)

- What should we learn?
 - p(r,s' | s,a)
 - $V_{\pi}(s)$
 - $q_{\pi}(s, a)$

Monte-Carlo RL

- Just like N+1 heuristic:
 - Get all trajectories containing particular (s, a)
 - Estimate $G_t(s, a)$ for each trajectory
 - Average them to get estimation of expectation



Monte-Carlo RL

- MC methods learn directly from episodes of experience
- MC is model-free: no knowledge of MDP transitions / rewards
- MC learns from complete episodes: no bootstrapping
- MC uses the simplest possible idea: value = mean return
- Note: can only apply MC to episodic MDPs
 - All episodes must terminate

Temporal Difference

- Just like in the 'incremental mean' example we can improve $q_{\pi}(s, \alpha)$ iteratively:

$$q_*(s, a) = \sum_{r,s'} p(r, s' \mid s, a) [r + \gamma \max_{a'} q_*(s', a')]$$

• We don't have $p(r, s' \mid s, a)$ to compute 'fair' expectation, so what should we do?

Temporal Difference

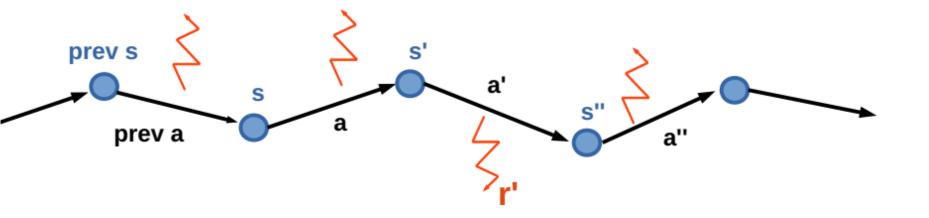
$$\sum_{r,s'} p(r,s' \mid s,a)[r + \gamma \max_{a'} q_*(s',a')] \approx$$

$$\approx \frac{1}{N} \sum_{i} r_{i} + \gamma \max_{a'} Q(s'_{i}, a')$$

 One more trick: use alpha-smoothing for updating Q-values.

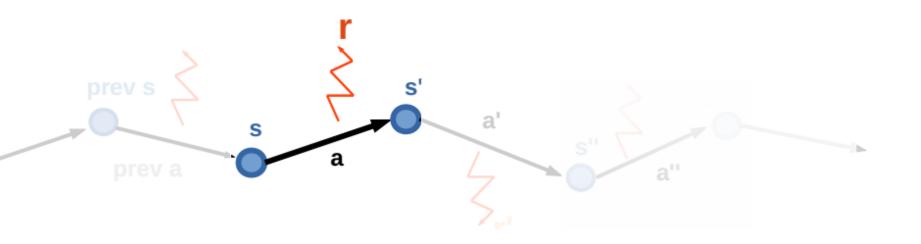
$$Q(s_t, a_t) = \alpha (r_t + \gamma \max_{a'} Q(s_{t+1}, a')) + (1 - \alpha) Q(s_t, at)$$

Q-learning



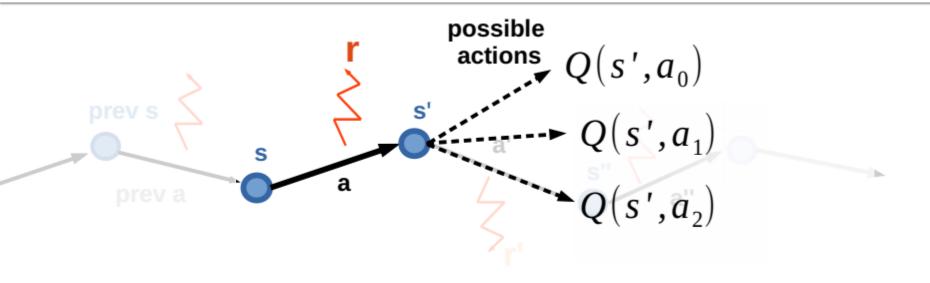
- Works on a sequence of
 - states (s)
 - actions (a)
 - rewards (r)

Q-learning



- Initialize Q(s, a) with zeros
- Cycle:
 - Sample <s, a, r, s'> from environment

Q-learning



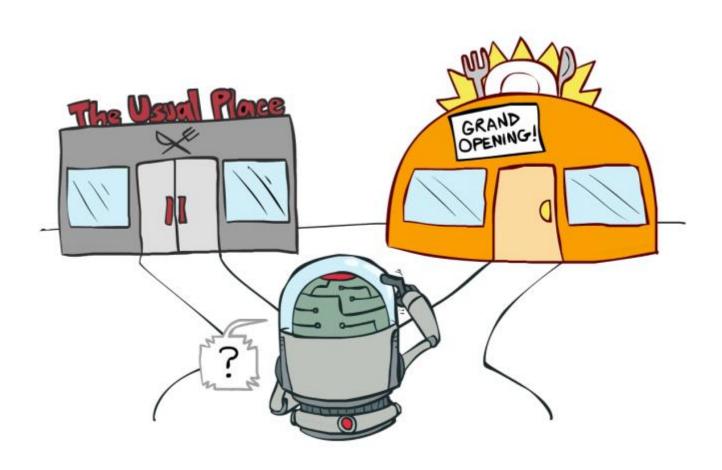
- Initialize Q(s, a) with zeros
- Cycle:
 - Sample <s, a, r, s'> from environment
 - Compute $\widehat{Q}(s,a) = r(s,a) + \gamma \max_{a_i} Q(s',ai)$
 - Update: $Q(s_t, at) = \alpha \hat{Q}(s, a) + (1 \alpha) Q(s_t, at)$

MC vs TD

- TD can learn before knowing the final outcome
 - TD can learn online after every step
 - MC must wait until end of episode before return is known
- TD can learn without the final outcome
 - TD can learn from incomplete sequences
 - MC can only learn from complete sequences
 - TD works in continuing (non-terminating) environments
 - MC only works for episodic (terminating) environments

Exploration/Exploitation Revisited

 Balance between using what you learned and trying to find something even better



Exploration/Exploitation Revisited

Strategies:

- ε-greedy
 With probability ε take random action, otherwise take optimal action.
- ε- dithering
 Adding random noise to Q-values with ε probability

Exploration/Exploitation over time

 If you want to converge to optimal policy you need to gradually reduce exploration.

Example:

Initialize ε -greedy ε = 0.5, then gradually reduce it

- If $\varepsilon \rightarrow o$, it's **greedy in the limit**
- Be careful with non-stationary environments

Temporal Difference Learning

- TD methods learn directly from episodes of experience
- TD is model-free: no knowledge of MDP transitions / rewards
- TD learns from incomplete episodes, by bootstrapping
- TD updates a guess towards a guess

Reinforcement Learning in the Wild

- Reinforcement learning can be used to solve large problems, e.g.
 - Backgammon: 1020 states
 - Computer Go: 10170 states
 - Helicopter: continuous state space
- How can we scale up the model-free methods for prediction and control?

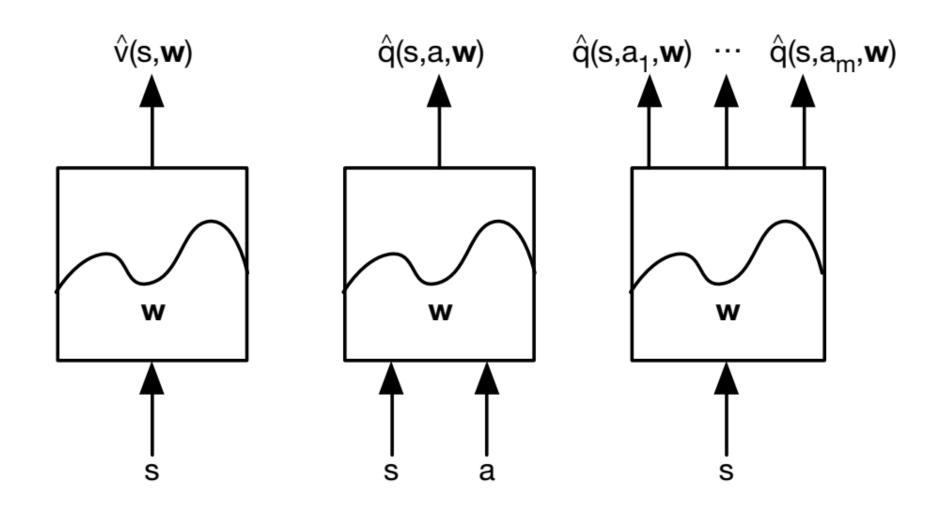
Curse of dimensionality in RL

<u>Problem:</u>

- State space is usually large, sometimes continuous.
- How about action space?
- However, states do have a structure, similar states have similar action outcomes

What should we do?

Types of Value Function Approximation



Which class of function to choose?

- There are many function approximators, e.g.
 - Linear combinations of features
 - Neural network
 - Decision tree
 - Nearest neighbor
 - Fourier / wavelet bases
 - •

Gradient Descent

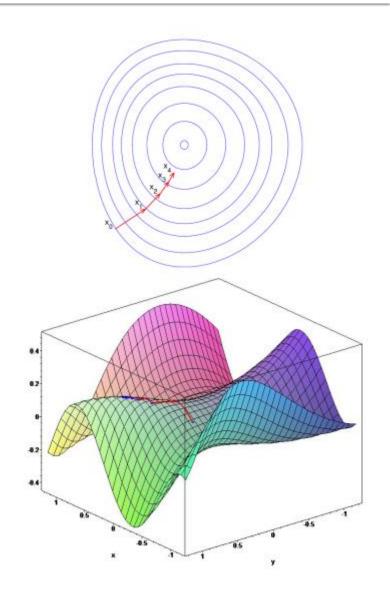
- Let J(w) be a differentiable function of parameter vector w
- Define the gradient of J(w) to be

$$\nabla_{\mathbf{w}} J(\mathbf{w}) = \begin{pmatrix} \frac{\partial J(\mathbf{w})}{\partial \mathbf{w}_1} \\ \vdots \\ \frac{\partial J(\mathbf{w})}{\partial \mathbf{w}_n} \end{pmatrix}$$

- To find a local minimum of J(w):
 - Adjust w in direction of -ve gradient

$$\Delta \mathbf{w} = -\frac{1}{2} \alpha \nabla_{\mathbf{w}} J(\mathbf{w})$$

where α is a step-size parameter



SGD for Value Function approximation

• Goal: find parameter vector **w** minimizing mean-squared error between approximate value $v'(s, \mathbf{w})$ and true value $v_{\pi}(s)$

$$J(\mathbf{w}) = \mathbb{E}_{\pi} \left[(v_{\pi}(S) - \hat{v}(S, \mathbf{w}))^2 \right]$$

Gradient descent finds a local minimum

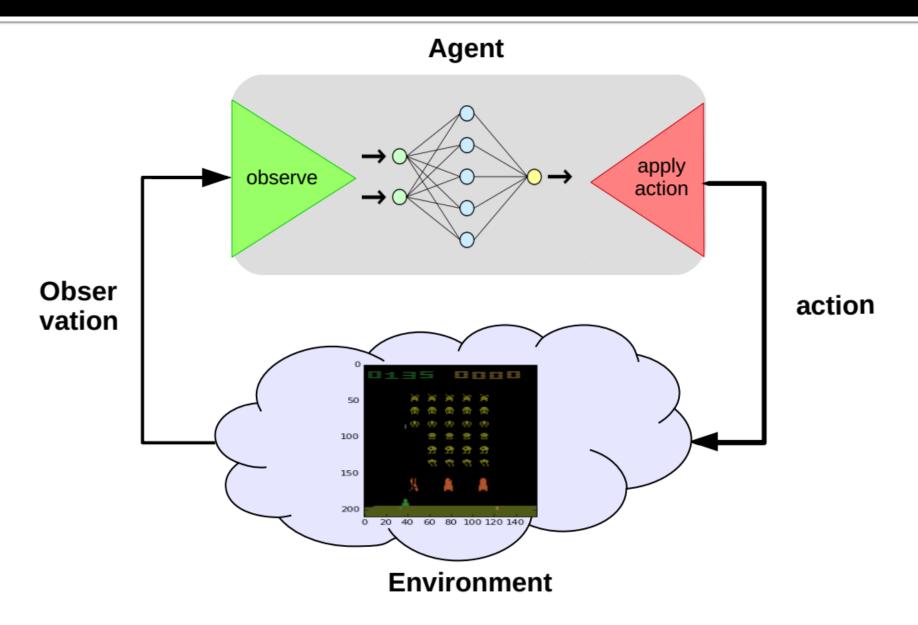
$$\Delta \mathbf{w} = -\frac{1}{2} \alpha \nabla_{\mathbf{w}} J(\mathbf{w})$$
$$= \alpha \mathbb{E}_{\pi} \left[(v_{\pi}(S) - \hat{v}(S, \mathbf{w})) \nabla_{\mathbf{w}} \hat{v}(S, \mathbf{w}) \right]$$

- Stochastic gradient descent *samples* the gradient

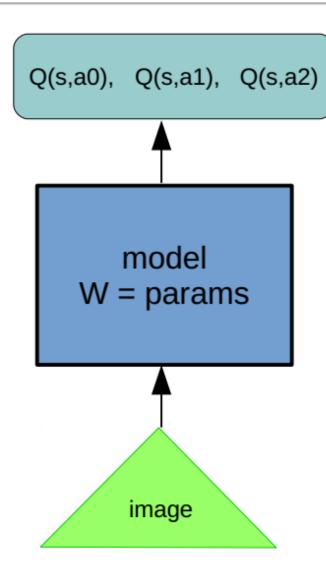
$$\Delta \mathbf{w} = \alpha(\mathbf{v}_{\pi}(S) - \hat{\mathbf{v}}(S, \mathbf{w})) \nabla_{\mathbf{w}} \hat{\mathbf{v}}(S, \mathbf{w})$$

Expected update is equal to full gradient update

Atari again



Approximate Q-learning



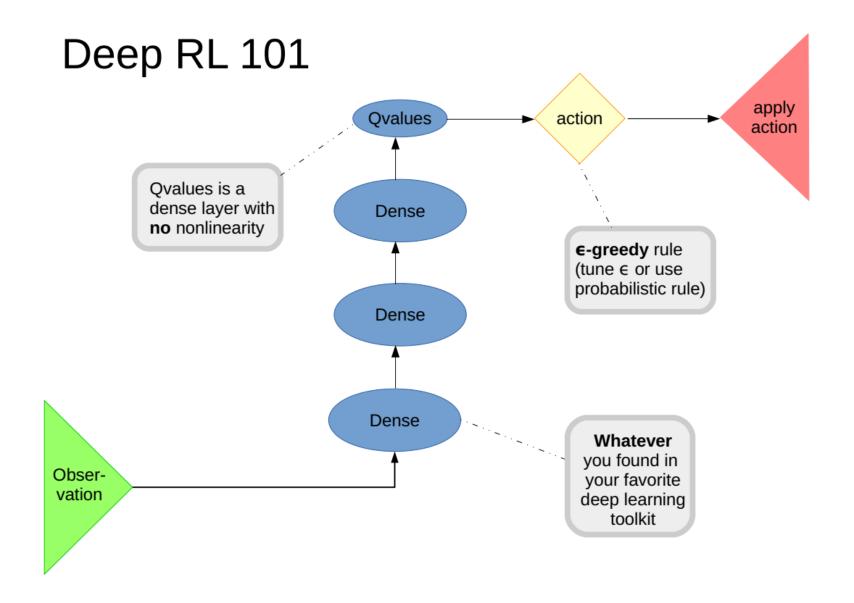
- Initialize W.
- Cycle:
 - Sample <s, a, r, s'> from environment
 - Compute $\widehat{m{Q}}(s,a) = r(s,a) + \gamma \max_{a_i} m{Q}(s',ai)$
 - Objective:

$$L = [Q(s_t, at) - \widehat{Q}(s_t, at)]^2$$

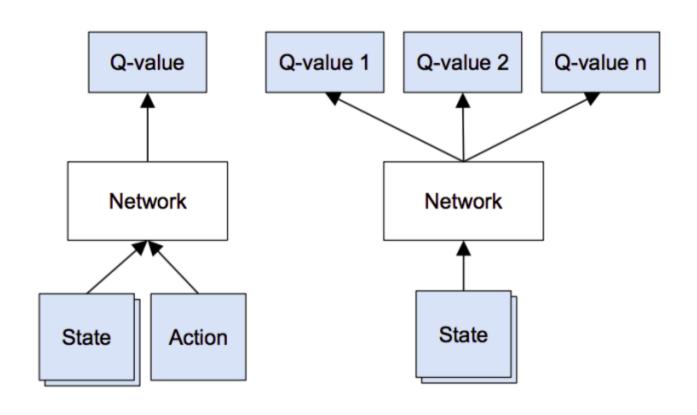
SGD Update:

$$W_{t+1} = W_t - \alpha \frac{\partial L}{\partial wt}$$

RL Mechanics

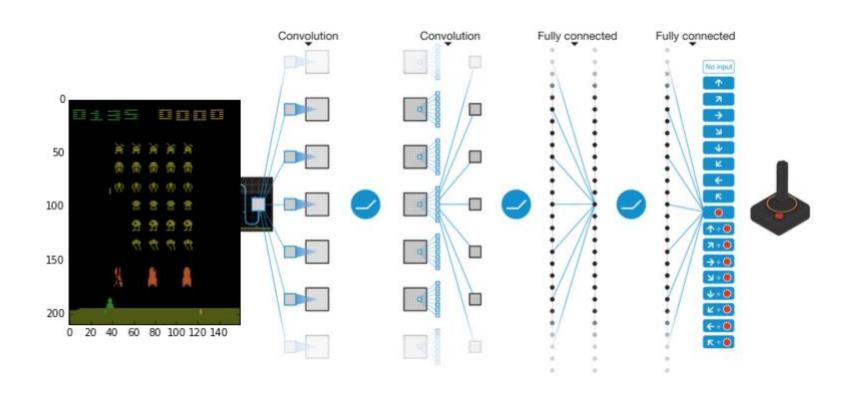


Architectures



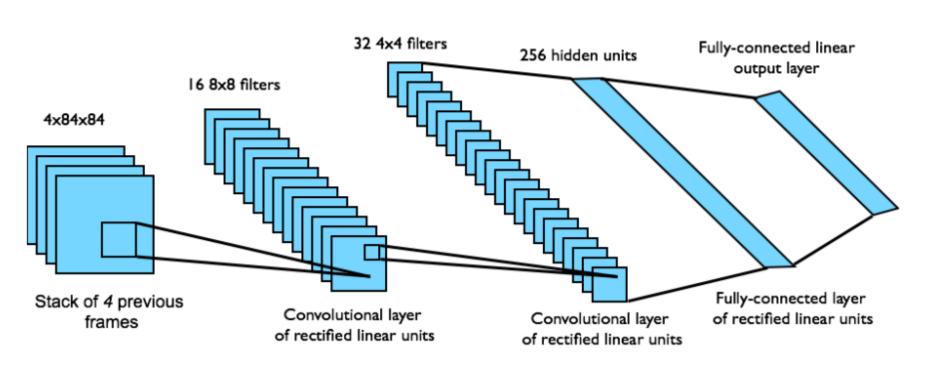
Given **(s,a)** Predict Q(s,a) Given **s** predict all q-values Q(s,a0), Q(s,a1), Q(s,a2)

From theory to practice: DQN case

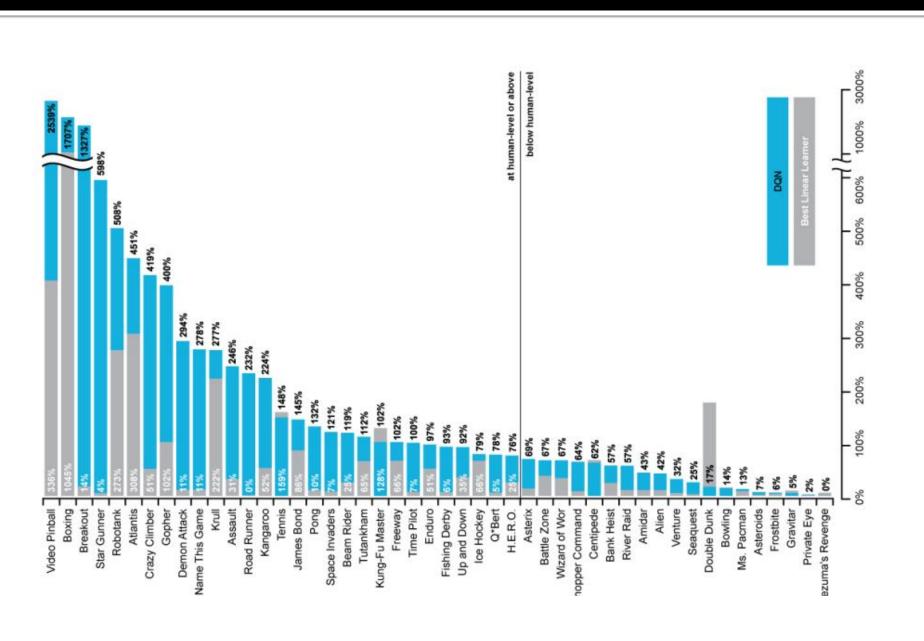


DQN: Atari

- End-to-end learning of values Q(s, a) from pixels s
- Input state s is stack of raw pixels from last 4 frames
- Output is Q(s, a) for **18** joystick/button positions
- Reward is change in score for that step



DQN results on Atari



DQN: under the hood

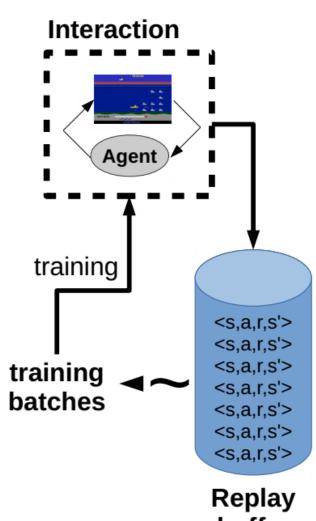
- DQN uses experience replay and fixed Q-targets:
 - Take action a_t according to e-greedy policy
 - Store transition $(s_t, \alpha_t, r_{t+1}, s_{t+1})$ in replay memory D
 - Sample random mini-batch of transitions (s, α, r, s') from D
 - Compute Q-learning targets w.r.t. old, fixed parameters w-
 - Optimize MSE between Q-network and Q-learning targets

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

Using variant of stochastic gradient descent

Experience Replay

- Idea: store several past interactions <*s*,*α*,*r*,*s*'>
- Train on random subsamples
- Any +/- ?



buffer

DQN: Atari Breakout

https://www.youtube.com/watch?v=TmPfTpjtdgg