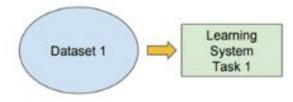
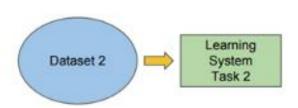
Reinforcement Learning Hands-On



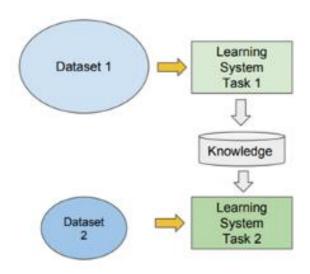
Traditional ML vs Transfer Learning

- Isolated, single task learning:
 - Knowledge is not retained or accumulated. Learning is performed w.o. considering past learned knowledge in other tasks





- Learning of a new tasks relies on the previous learned tasks:
 - Learning process can be faster, more accurate and/or need less training data



What to transfer

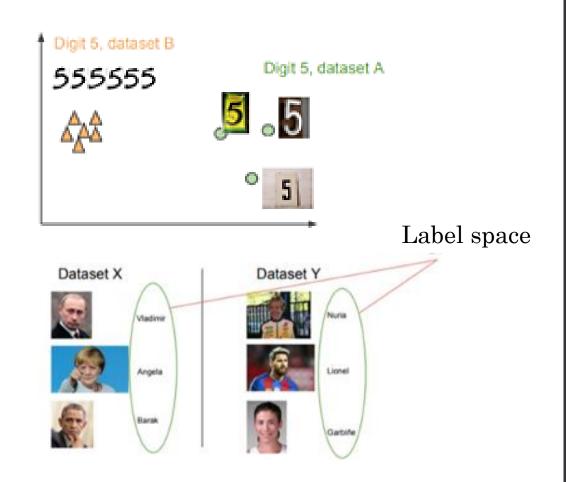
- Source-specific
- Common domain source-target

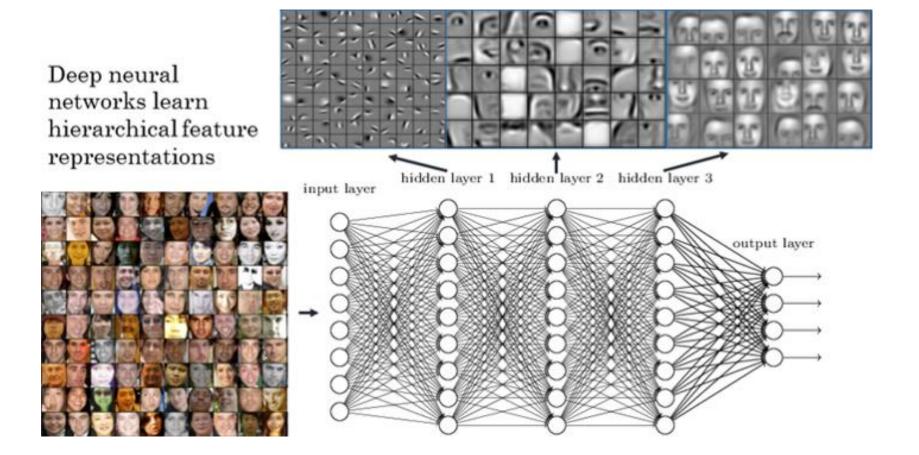
When to transfer

Negative transfer

How to transfer

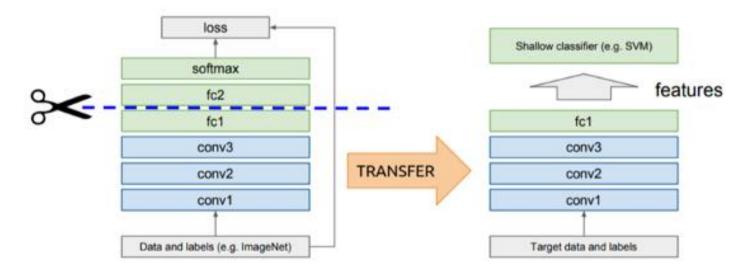
- Freeze Layers
- Fine-tuning Layers





Idea: use outputs of one or more layers of a network trained on a different task as generic feature detectors. Train a new shallow model on these features.

Assumes that $D_S = D_T$



Freeze or fine-tune?

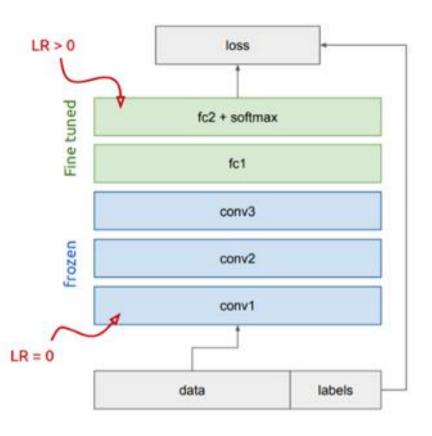
Bottom n layers can be frozen or fine tuned.

- Frozen: not updated during backprop
- Fine-tuned: updated during backprop

Which to do depends on target task:

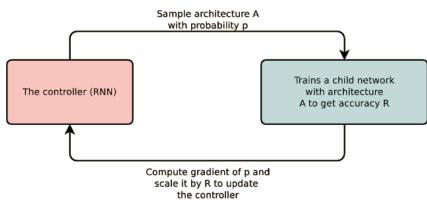
- Freeze: target task labels are scarce, and we want to avoid overfitting
- · Fine-tune: target task labels are more plentiful

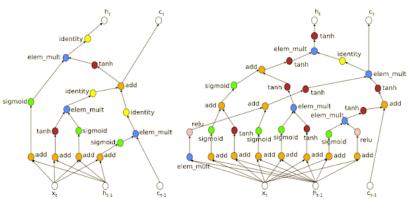
In general, we can set learning rates to be different for each layer to find a tradeoff between freezing and fine tuning



NASNet and AutoML

AI build other AI Using Reinforcement Learning

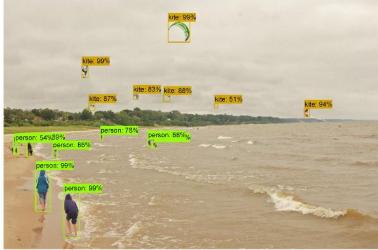


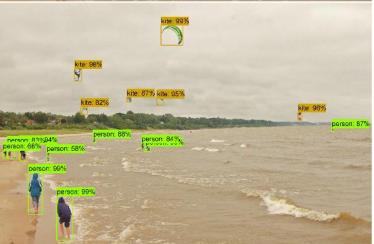


 $Learning\ Transferable\ Architectures\ for\ Scalable\ Image\ Recognition,\ Jul\ 2017\ Google, \\ \underline{https://arxiv.org/pdf/1707.07012.pdf}$

https://arxiv.org/pdf/1707.07012.pdf https://research.googleblog.com/2017/05/using-machine-learning-to-explore.html

Neural Architecture Search





Pipeline:

Faster R-CNN Feature map Generator:

Inception-ResNet

Pipeline:Faster R-CNN

Feature map Generator:

NASNet



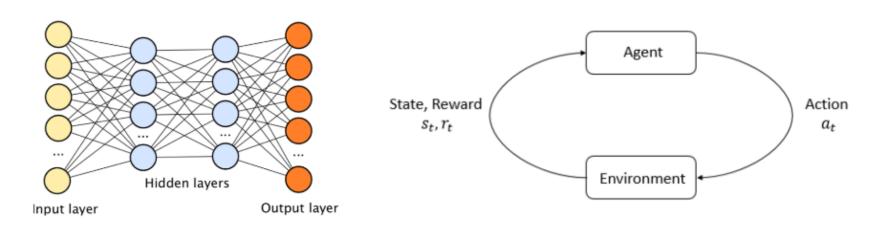


https://blog.openai.com/openai-charter/

What is Deep RL?

Deep RL is the combination of reinforcement learning (RL) with deep learning

- Reinforcement learning is about solving problems by trial and error
- Deep learning is about using deep neural networks to solve problems
- **Deep reinforcement learning** trains deep neural networks with trial and error



Deep neural network¹ and RL interaction loop

What is Deep RL?

RL is useful when

- you have a sequential decision-making problem
- you do not know the optimal behavior already¹
- but you can still evaluate whether behaviors are good or bad

Deep learning is useful when

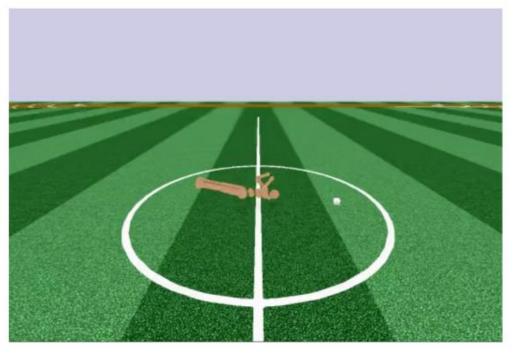
- you want to approximate a function
- function requires "intelligence"
- inputs and/or outputs are high-dimensional
- lots of data is available

Uses Deep RL ...

Deep RL can...

- Play video games from raw pixels
- Control robots in simulation and in the real world
- Play Go, Dota, and Starcraft at superhuman levels





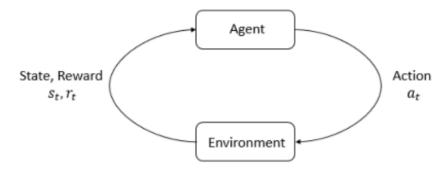
RL Concepts

What are the core parts for understand RL:

- Observations and actions spaces
- Policies
- Trajectories
- Reward and return
- RL optimization problem
- Value and Action-Value Functions

RL Setup

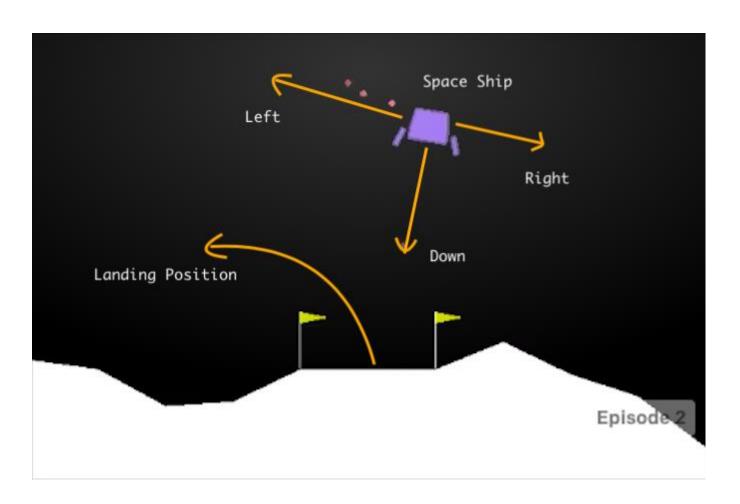
An agent interacts with an environment.



```
obs = env.reset()
done = False
while not(done):
   act = agent.get_action(obs)
   next_obs, reward, done, info = env.step(act)
   obs = next_obs
```

- The goal of the agent is to maximize cumulative reward (called return).
- The agent figures out how to attain its goal by trial and error.
- Reinforcement learning (RL) is a field of study for algorithms that do that.

Environment



Get environment

- pip install gym
- conda install pystan
- conda install git
- pip install git+https://github.com/Kojoley/atari-py.git
- conda install swig
- pip install box2d-py && pip install Box2D
- pip install gym[all]
- pip install pyglet==1.2.4
- pip install gym[box2d]

Policy

A **policy** π is a rule for selecting actions. It can be either

- **stochastic**, which means that it gives a probability distribution over actions, and actions are selected randomly based on that distribution $(a_t \sim \pi(\cdot|s_t))$,
- or **deterministic**, which means that π directly maps to an action $(a_t = \pi(s_t))$.

Examples of policies:

Stochastic policy over discrete actions:

```
obs = tf.placeholder(shape=(None, obs_dim), dtype=tf.float32)
net = mlp(obs, hidden_dims=(64,64), activation=tf.tanh)
logits = tf.layers.dense(net, units=num_actions, activation=None)
actions = tf.squeeze(tf.multinomial(logits=logits,num_samples=1), axis=1)
```

Deterministic policy for a vector-valued continuous action:

```
obs = tf.placeholder(shape=(None, obs_dim), dtype=tf.float32)
net = mlp(obs, hidden_dims=(64,64), activation=tf.tanh)
actions = tf.layers.dense(net, units=act_dim, activation=None)
```

Trajectories

• A trajectory τ is a sequence of states and actions in an environment:

$$\tau = (s_0, a_0, s_1, a_1, ...).$$

• The initial state s_0 is sampled from a start state distribution μ :

$$s_0 \sim \mu(\cdot)$$
.

 State transitions depend only on the most recent state and action. They could be deterministic:

$$s_{t+1} = f(s_t, a_t),$$

or stochastic:

$$s_{t+1} \sim P(\cdot|s_t, a_t).$$

A trajectory is sometimes also called an episode or rollout.

Reward

The **reward** function of an environment measures how good state-action pairs are:

$$r_t = R(s_t, a_t).$$

The **return** of a trajectory is a measure of cumulative reward along it. There are two main ways to compute return:

• Finite horizon undiscounted sum of rewards::

$$R(\tau) = \sum_{t=0}^{T} r_t$$

Infinite horizon discounted sum of rewards:

$$R(\tau) = \sum_{t=0}^{\infty} \gamma^t r_t$$

where $\gamma \in (0,1)$. This makes rewards less valuable if they are further in the future. (Why would we ever want this? Think about cash: it's valuable to have it sooner rather than later!)

Reward-to-go

A closely-related quantity is **reward-to-go**, which is "return starting from a state or timestep":

$$R_t = \sum_{t'=t}^T r_{t'}$$

or

$$R_t = \sum_{t'=t}^{\infty} \gamma^{t'-t} r_{t'}$$

What to learn?

In RL we want a policy which maximizes expected return. Thus the performance measure is

$$J(\pi) = \mathop{\rm E}_{\tau \sim \pi} \left[R(\tau) \right],$$

and the **optimal policy** π^* is:

$$\pi^* = \arg\max_{\pi} J(\pi)$$

Note that by $\tau \sim \pi$, we mean

$$s_0 \sim \mu(\cdot), \quad a_t \sim \pi(\cdot|s_t), \quad s_{t+1} \sim P(\cdot|s_t, a_t).$$

V^{π} function and Q^{π} function

Following a policy produces sample trajectories (or paths) s_0 , a_0 , r_0 , s_1 , a_1 , r_1 , ...

How good is a state?

The **value function** at state s, is the expected cumulative reward from following the policy from state s:

 $V^{\pi}(s) = \mathbb{E}\left[\sum_{t \geq 0} \gamma^t r_t | s_0 = s, \pi
ight]$

How good is a state-action pair?

The **Q-value function** at state s and action a, is the expected cumulative reward from taking action a in state s and then following the policy:

$$Q^{\pi}(s,a) = \mathbb{E}\left[\sum_{t \geq 0} \gamma^t r_t | s_0 = s, a_0 = a, \pi
ight]$$

Advantage function

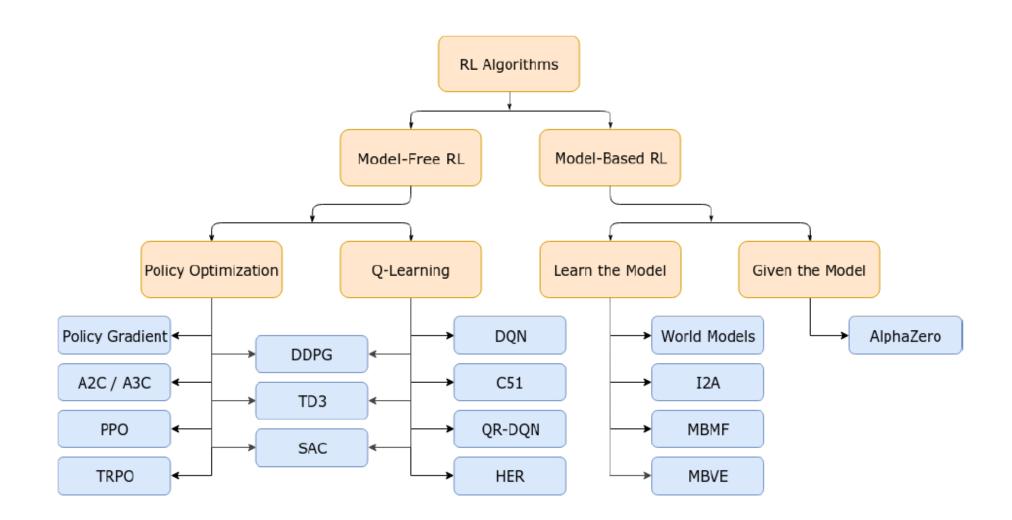
Value and action-value functions are connected:

$$V^{\pi}(s) = \mathop{\mathrm{E}}_{a \sim \pi} \left[Q^{\pi}(s, a) \right]$$

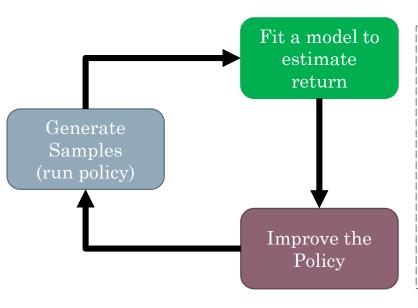
The advantage function for a policy tells you how much better or worse one action is than average:

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

Deep RL Tree



RL flow



Policy Optimization

• Policy $\pi_{\theta}(a|s)$

- Optimize θ using objective function $I(\pi_{\theta})$
- On-Policy, only use data collected from recent policy
- Can involve learning a value function $V^{\pi}(s)$
- A2C/A3C , TRPO

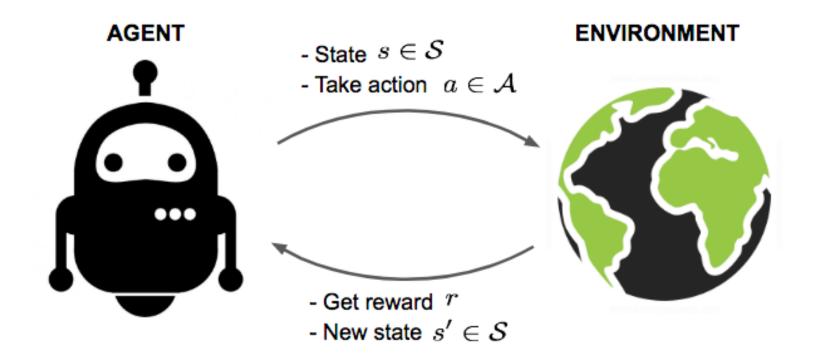
- Q-Learning
- Learn to optimize an action-value function $Q_{\theta}(s, a)$
- Function based on *Bellman-equation*
- Off-policy as each update use data collected at any point
- · DQN, Double-DQN

$$J(\pi_{\theta}) = \sum_{t=1}^{T} \gamma^t r_t$$

$$\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\pi_{\theta})$$

$$fit \ Q_{\pi_{\theta}}(s,a|\theta)$$

$$\pi_{\theta}(s) = argmax(Q_{\pi_{\theta}}(s, a))$$



DQN

Q-Learning

- Run policy: Step in env with action from Q_{θ} , store to replay buffer
- Evaluate policy: Update Q_{θ} to minimize Bellman error, using all previous data (off-policy)
- Improve policy: $a^* = \arg \max_a Q_{\theta}(s, a)$

Q-learning: Use a function approximator to estimate the action-value function

$$Q(s,a;\theta) pprox Q^*(s,a)$$
 function parameters (weights)

If the function approximator is a deep neural network => deep q-learning!

DQN

- Collect experience in the environment using a policy which trades off between acting randomly and acting according to current Q_{θ}
- Interleave data collection with updates to Q_{θ} to minimize Bellman error by bootstrapping:

$$\min_{\theta} \sum_{(s,a,s',r)\in\mathcal{D}} \left(Q_{\theta}(s,a) - y(s',r)\right)^{2}$$

where

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

and gradients don't propagate through y.

Experience replay:

- Data distribution changes over time: as your Q function gets better and you exploit this, you visit different (s, a, s', r) transitions than you did earlier
- Stabilize learning by keeping old transitions in a replay buffer, and taking minibatch gradient descent on mix of old and new transitions

Target networks:

- Bootstrapping with function approximators is unstable!
- It's like regression, but it's not: targets depend on parameters θ —so an update to Q changes the target!

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

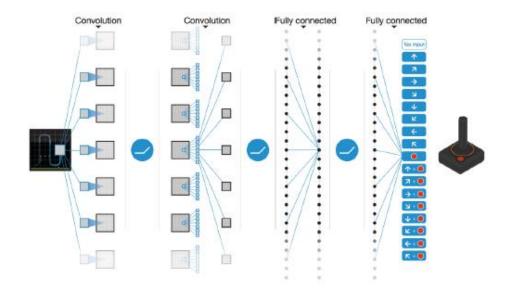
- Stabilize it by holding the target fixed for a while: keep a separate target network, $Q_{\theta_{targ}}$, and every k steps update $\theta_{targ} \leftarrow \theta$
- If unstable, why do it? Because it works well! (Plus theory, later)

What we've described so far requires the computation of

$$\max_{a} Q_{\theta}(s, a),$$

both for selecting actions and calculating update targets.

Problem: for continuous action spaces, this is nontrivially hard! DQN therefore only applies for **discrete actions**, because we can output one Q-value per action from the network.



Algorithm 2 Deep Q-Learning

```
Randomly generate Q-function parameters \theta

Set target Q-network parameters \theta_{targ} \leftarrow \theta

Make empty replay buffer \mathcal{D}

Receive observation s_0 from environment

for t=0,1,2,... do

With probability \epsilon, select random action a_t; otherwise select a_t = \arg\max_a Q_\theta(s_t,a)

Step environment to get s_{t+1}, r_t and end-of-episode signal d_t

Linearly decay \epsilon until it reaches final value \epsilon_f

Store (s_t, a_t, r_t, s_{t+1}, d_t) \rightarrow \mathcal{D}

Sample mini-batch of transitions B = \{(s, a, r, s', d)_i\} from \mathcal{D}

For each transition in B, compute
y = \begin{cases} r & \text{transition is terminal } (d = \text{True}) \\ r + \gamma \max_{a'} Q_{\theta_{targ}}(s', a') & \text{otherwise} \end{cases}
```

Update Q by gradient descent on regression loss:

$$\theta \leftarrow \theta - \alpha \nabla_{\theta} \sum_{(s,a,y) \in B} (Q_{\theta}(s,a) - y)^2$$

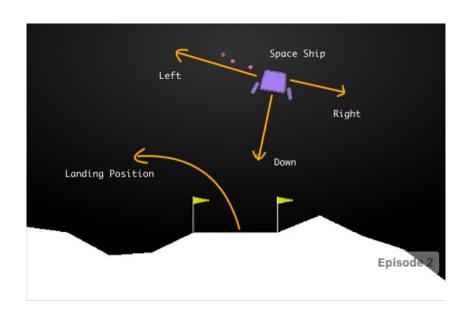
```
if t\%t_{update} = 0 then

Set \theta_{targ} \leftarrow \theta

end if

end for
```

Example



• Fuel is infinite, so an agent can learn to fly and then land on its first attempt.

Observation

- Landing pad is always at coordinates (0,0).
- Coordinates are the first two numbers in state vector.

Reward

- From the top of the screen to landing pad and zero speed is about 100..140 points.
- If lander moves away from landing pad it loses reward back.
- Episode finishes if the lander crashes or comes to rest, receiving additional -100 or +100 points.
- Each leg ground contact is +10.
- Firing main engine is -0.3 points each frame.
- Solved is 200 points.

Actions

• Four discrete actions available: do nothing, fire left orientation engine, fire main engine, fire right orientation engine.

Algorithm 2 Deep Q-Learning

Randomly generate Q-function parameters θ Set target Q-network parameters $\theta_{targ} \leftarrow \theta$ Make empty replay buffer \mathcal{D} Receive observation s_0 from environment



```
class DQN:
  def __init__(self, action_space, state_space):
      self.action space = action space
      self.state space = state space
      self.epsilon = 1.0
     self.gamma = .99
     self.batch size = 64
     self.epsilon min = .01
     self.lr = 0.001
     self.epsilon_decay = .996
     self.memory = deque(maxlen=1000000)
     self.model = self.build model()
   def build model(self):
      model = Sequential()
      model.add(Dense(150, input_dim=self.state_space, activation=relu))
      model.add(Dense(120, activation=relu))
      model.add(Dense(self.action_space, activation=linear))
      model.compile(loss='mse', optimizer=adam(lr=self.lr))
      return-model
```

$\overline{\mathrm{DQN}}$

```
for t=0,1,2,... do With probability \epsilon, select random action a_t; otherwise select a_t=\arg\max_a Q_{\theta}(s_t,a) Step environment to get s_{t+1}, r_t and end-of-episode signal d_t
```

```
def act(self, state):
    if np.random.rand() <= self.epsilon:
        return random.randrange(self.action_space)
        act_values = self.model.predict(state)
        return np.argmax(act_values[0])</pre>
```

```
agent = DQN(env.action_space.n, env.observation_space.shape[0])
for e in range(episode):
    state = env.reset()
    state = np.reshape(state, (1, 8))
    score = 0
    max_steps = 3000
    for i in range(max_steps):
        action = agent.act(state)
        env.render()
        next_state, reward, done, _ = env.step(action)
        score += reward
        next_state = np.reshape(next_state, (1, 8))
```

```
for t=0,1,2,... do

With probability \epsilon, select random action a_t; otherwise select a_t=\arg\max_a Q_\theta(s_t,a)

Step environment to get s_{t+1}, r_t and end-of-episode signal d_t

Linearly decay \epsilon until it reaches final value \epsilon_f

Store (s_t, a_t, r_t, s_{t+1}, d_t) \to \mathcal{D}

if self.epsilon > self.epsilon_min:

self.epsilon *= self.epsilon_decay
```

```
def remember(self, state, action, reward, next_state, done):
    self.memory.append((state, action, reward, next_state, done))
```

agent.remember(state, action, reward, next_state, done)



Sample mini-batch of transitions $B = \{(s, a, r, s', d)_i\}$ from \mathcal{D} For each transition in B, compute

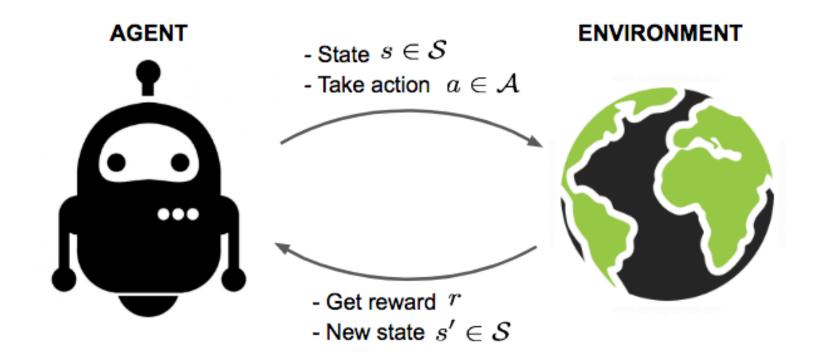
$$y = \begin{cases} r & \text{transition is terminal } (d = \text{True}) \\ r + \gamma \max_{a'} Q_{\theta_{targ}}(s', a') & \text{otherwise} \end{cases}$$

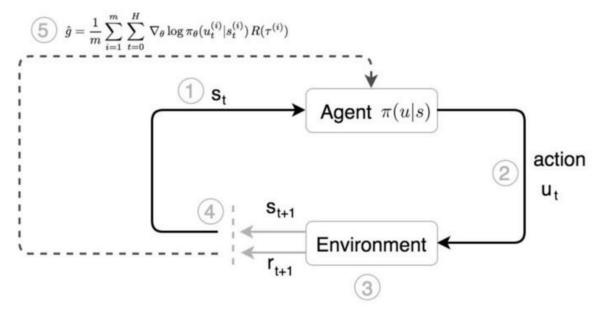


Update Q by gradient descent on regression loss:

$$\theta \leftarrow \theta - \alpha \nabla_{\theta} \sum_{(s,a,y) \in B} (Q_{\theta}(s,a) - y)^2$$

```
def replay(self):
  if len(self.memory) < self.batch_size:</pre>
      return
  minibatch = random.sample(self.memory, self.batch size)
  states = np.array([i[0] for i in minibatch])
  actions = np.array([i[1] for i in minibatch])
  rewards = np.array([i[2] for i in minibatch])
  next_states = np.array([i[3] for i in minibatch])
  dones = np.array([i[4] for i in minibatch])
  states = np.squeeze(states)
  next_states = np.squeeze(next_states)
  q_val = (np.amax(self.model.predict_on_batch(next_states), axis=1))
  targets = rewards + self.gamma*q val*(1-dones)
  targets_full = self.model.predict_on_batch(states)
  ind = np.array([i for i in range(self.batch_size)])
  targets full[[ind], [actions]] = targets
  self.model.fit(states, targets full, epochs=1, verbose=0)
```





Policy Optimization

- Run policy: Collect trajectories $\tau \sim \pi_{\theta}$
- Evaluate policy: Estimate $V^{\pi_{\theta}}$, $A^{\pi_{\theta}}$ using current trajectories (on-policy)
- Improve policy: Increase likelihood of actions with high advantage

Goal: derive an expression for $\nabla_{\theta} J(\pi_{\theta})$ which we can compute with a sample estimate, as the basis for a direct gradient ascent algorithm

$$\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\pi_{\theta})$$

Well what happens if we just...

$$\nabla_{\theta} J(\pi_{\theta}) = \nabla_{\theta} \mathop{\to}_{\tau \sim \pi_{\theta}} [R(\tau)]$$

Problem: parameters are in distribution!

Solution: expand expectation into integral, use log-derivative trick

$$\nabla_{\theta} J(\pi_{\theta}) = \nabla_{\theta} \mathop{\to}_{\tau \sim \pi_{\theta}} [R(\tau)]$$

$$= \nabla_{\theta} \int d\tau P(\tau | \pi_{\theta}) R(\tau)$$

$$= \int d\tau \nabla_{\theta} P(\tau | \pi_{\theta}) R(\tau)$$

$$= \int d\tau P(\tau | \pi_{\theta}) \nabla_{\theta} \log P(\tau | \pi_{\theta}) R(\tau)$$

$$= \mathop{\to}_{\tau \sim \pi_{\theta}} [\nabla_{\theta} \log P(\tau | \pi_{\theta}) R(\tau)]$$

But are we done yet? No! Still need to compute $\nabla_{\theta} \log P(\tau | \pi_{\theta})$

What is $P(\tau|\pi_{\theta})$?

$$P(\tau|\pi_{\theta}) = \mu(s_0) \prod_{t=0}^{T} P(s_{t+1}|s_t, a_t) \pi_{\theta}(a_t|s_t)$$

Thus:

$$\nabla_{\theta} \log P(\tau | \pi_{\theta}) = \nabla_{\theta} \log \left(\mu(s_{0}) \prod_{t=0}^{T} P(s_{t+1} | s_{t}, a_{t}) \pi_{\theta}(a_{t} | s_{t}) \right)$$

$$= \nabla_{\theta} \left(\log \mu(s_{0}) + \sum_{t=0}^{T} \left(\log P(s_{t+1} | s_{t}, a_{t}) + \log \pi_{\theta}(a_{t} | s_{t}) \right) \right)$$

$$= \nabla_{\theta} \log \mu(s_{0}) + \sum_{t=0}^{T} \left(\nabla_{\theta} \log P(s_{t+1} | s_{t}, a_{t}) + \nabla_{\theta} \log \pi_{\theta}(a_{t} | s_{t}) \right)$$

$$= \nabla_{\theta} \log \mu(s_{0}) + \sum_{t=0}^{T} \left(\nabla_{\theta} \log P(s_{t+1} | s_{t}, a_{t}) + \nabla_{\theta} \log \pi_{\theta}(a_{t} | s_{t}) \right)$$

$$\therefore \nabla_{\theta} \log P(\tau | \pi_{\theta}) = \sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t} | s_{t})$$

Putting it all together so far:

$$\nabla_{\theta} J(\pi_{\theta}) = \mathop{\mathbf{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) R(\tau) \right]$$

So we could estimate with:

$$abla_{ heta} J(\pi_{ heta}) pprox rac{1}{|\mathcal{D}|} \sum_{ au \in \mathcal{D}} \sum_{t=0}^{T}
abla_{ heta} \log \pi_{ heta}(a_{t}|s_{t}) R(au)$$

But not good enough! Variance will be high.

Insight: future actions and past rewards should be uncorrelated. That is:

for
$$t > t'$$
, $\mathrm{E}\left[\nabla_{\theta} \log \pi_{\theta}(a_t|s_t)r_{t'}\right] = 0$

$$\therefore \nabla_{\theta} J(\pi_{\theta}) = \mathop{\mathbb{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \sum_{t'=t}^{T} r_{t'} \right]$$

REINFORCE algorithm:



- 1. sample $\{\tau^i\}$ from $\pi_{\theta}(\mathbf{a}_t|\mathbf{s}_t)$ (run the policy)
- 2. $\nabla_{\theta} J(\theta) \approx \sum_{i} \left(\sum_{t} \nabla_{\theta} \log \pi_{\theta}(\mathbf{a}_{t}^{i} | \mathbf{s}_{t}^{i}) \right) \left(\sum_{t} r(\mathbf{s}_{t}^{i}, \mathbf{a}_{t}^{i}) \right)$
- 3. $\theta \leftarrow \theta + \alpha \nabla_{\theta} J(\theta)$

Policy gradient:

```
# Given:
# actions - (N*T) x Da tensor of actions
# states - (N*T) x Ds tensor of states
# q_values - (N*T) x 1 tensor of estimated state-action values
# Build the graph:
logits = policy.predictions(states) # This should return (N*T) x Da tensor of action logits
negative_likelihoods = tf.nn.softmax_cross_entropy_with_logits(labels=actions, logits=logits)
weighted_negative_likelihoods = tf.multiply(negative_likelihoods, q_values)
loss = tf.reduce_mean(weighted_negative_likelihoods)
gradients = loss.gradients(loss, variables)
```

What we have currently: "Reward-to-Go" policy gradient:

$$\nabla_{\theta} J(\pi_{\theta}) = \mathop{\mathbb{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \sum_{t'=t}^{T} r_{t'} \right]$$

Observe: expectation can be broken up, letting us transform "Reward-to-Go" into Q^{π} :

$$\nabla_{\theta} J(\pi_{\theta}) = \underset{\tau \sim \pi_{\theta}}{\mathbb{E}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \sum_{t'=t}^{T} r_{t'} \right]$$

$$= \sum_{t=0}^{T} \underset{\tau \sim \pi_{\theta}}{\mathbb{E}} \left[\nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \sum_{t'=t}^{T} r_{t'} \right]$$

$$= \sum_{t=0}^{T} \underset{\tau_{0:t} \sim \pi_{\theta}}{\mathbb{E}} \left[\nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \underset{\tau_{(t+1):T} \sim \pi_{\theta}}{\mathbb{E}} \left[\sum_{t'=t}^{T} r_{t'} \right] \right]$$

$$= \sum_{t=0}^{T} \underset{\tau_{0:t} \sim \pi_{\theta}}{\mathbb{E}} \left[\nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) Q^{\pi_{\theta}}(s_{t}, a_{t}) \right]$$

$$\therefore \nabla_{\theta} J(\pi_{\theta}) = \underset{\tau \sim \pi_{\theta}}{\mathbb{E}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) Q^{\pi_{\theta}}(s_{t}, a_{t}) \right]$$

What is a **baseline**? A function $b(s_t)$ with

$$\nabla_{\theta} J(\pi_{\theta}) = \mathop{\mathrm{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \left(Q^{\pi_{\theta}}(s_{t}, a_{t}) - b(s_{t}) \right) \right]$$

Claim: this works for any b! Proof:

$$\begin{split} & \underset{a_{t} \sim \pi_{\theta}(\cdot|s_{t})}{\text{E}} \left[\nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t})b(s_{t}) \right] = \underset{a_{t} \sim \pi_{\theta}(\cdot|s_{t})}{\text{E}} \left[\nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \right] b(s_{t}) \\ & = \left(\int da \pi_{\theta}(a_{t}|s_{t}) \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \right) b(s_{t}) \\ & = \left(\int da \nabla_{\theta} \pi_{\theta}(a_{t}|s_{t}) \right) b(s_{t}) \\ & = \left(\nabla_{\theta} \int da \pi_{\theta}(a_{t}|s_{t}) \right) b(s_{t}) \\ & = \left(\nabla_{\theta} \Upsilon_{\theta} \Upsilon_{\theta}(a_{t}|s_{t}) \right) b(s_{t}) \\ & = \left(\nabla_{\theta} \Upsilon_{\theta} \Upsilon_{\theta}(a_{t}|s_{t}) \right) b(s_{t}) \\ & = 0 \end{split}$$

If we choose $b = V^{\pi}$, we get the **advantage form** of the policy gradient:

$$\nabla_{\theta} J(\pi_{\theta}) = \mathop{\mathbb{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \left(Q^{\pi_{\theta}}(s_{t}, a_{t}) - V^{\pi_{\theta}}(s_{t}) \right) \right]$$
$$= \mathop{\mathbb{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) A^{\pi_{\theta}}(s_{t}, a_{t}) \right]$$

Why do we want this? Better signal in sample estimate: removes "stuff that would have happened anyway" from Q^{π}

What we've shown so far:

$$\nabla_{\theta} J(\pi_{\theta}) = \underset{\tau \sim \pi_{\theta}}{\text{E}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) R(\tau) \right]$$

$$= \underset{\tau \sim \pi_{\theta}}{\text{E}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) \sum_{t'=t}^{T} r_{t'} \right]$$

$$= \underset{\tau \sim \pi_{\theta}}{\text{E}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) Q^{\pi_{\theta}}(s_{t}, a_{t}) \right]$$

$$= \underset{\tau \sim \pi_{\theta}}{\text{E}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) A^{\pi_{\theta}}(s_{t}, a_{t}) \right]$$

Which form do we use? Almost always the last one.

$$\nabla_{\theta} J(\pi_{\theta}) = \mathop{\mathrm{E}}_{\tau \sim \pi_{\theta}} \left[\sum_{t=0}^{T} \nabla_{\theta} \log \pi_{\theta}(a_{t}|s_{t}) A^{\pi_{\theta}}(s_{t}, a_{t}) \right]$$

- Pushes up the probabilities of "good" actions and pushes down probabilities of "bad" actions
- To estimate, must sample on-policy

- Generalized Advantage Estimation (GAE)
- Natural Policy Gradients (NPG)

Algorithm 1 Vanilla Policy Gradient Algorithm

- 1: Input: initial policy parameters θ_0 , initial value function parameters ϕ_0
- 2: for k = 0, 1, 2, ... do
- Collect set of trajectories $\mathcal{D}_k = \{\tau_i\}$ by running policy $\pi_k = \pi(\theta_k)$ in the environment.
- Compute rewards-to-go \hat{R}_t .
- Compute advantage estimates, \hat{A}_t (using any method of advantage estimation) based on the current value function V_{ϕ_k} .
- Estimate policy gradient as 6:

$$\hat{g}_k = rac{1}{|\mathcal{D}_k|} \sum_{ au \in \mathcal{D}_k} \sum_{t=0}^T |
abla_{ heta} \log \pi_{ heta}(a_t|s_t)|_{ heta_k} \hat{A}_t.$$

Compute policy update, either using standard gradient ascent, 7:

$$\theta_{k+1} = \theta_k + \alpha_k \hat{g}_k,$$



or via another gradient ascent algorithm like Adam.

8: Fit value function by regression on mean-squared error:

$$\phi_{k+1} = rg \min_{\phi} rac{1}{|\mathcal{D}_k|T} \sum_{ au \in \mathcal{D}_k} \sum_{t=0}^T \left(V_{\phi}(s_t) - \hat{R}_t
ight)^2,$$

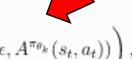
typically via some gradient descent algorithm.

9: end for

Proximal Policy Gradient

Algorithm 1 PPO-Clip

- 1: Input: initial policy parameters θ_0 , initial value function parameters ϕ_0
- 2: for k = 0, 1, 2, ... do
- Collect set of trajectories $D_k = \{\tau_i\}$ by running policy $\pi_k = \pi(\theta_k)$ in the environment.
- Compute rewards-to-go \bar{R}_t .
- Compute advantage estimates, \hat{A}_t (using any method of advantage estimation) based on the current value function V_{ϕ_k} .
- Update the policy by maximizing the PPO-Clip objective:



$$\theta_{k+1} = \arg\max_{\theta} \frac{1}{|\mathcal{D}_k|T} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^{T} \min\left(\frac{\pi_{\theta}(a_t|s_t)}{\pi_{\theta_k}(a_t|s_t)} A^{\pi_{\theta_k}}(s_t, a_t), \ g(\epsilon, A^{\pi_{\theta_k}}(s_t, a_t))\right),$$

typically via stochastic gradient ascent with Adam.

Fit value function by regression on mean-squared error:

$$\phi_{k+1} = \arg \min_{\phi} \frac{1}{|\mathcal{D}_k|T} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^{T} (V_{\phi}(s_t) - \hat{R}_t)^2,$$

typically via some gradient descent algorithm.

8: end for

Trust Region Policy Gradient

Algorithm 1 Trust Region Policy Optimization

- 1: Input: initial policy parameters θ_0 , initial value function parameters ϕ_0
- 2: Hyperparameters: KL-divergence limit δ , backtracking coefficient α , maximum number of backtracking steps K
- 3: for k = 0, 1, 2, ... do
- Collect set of trajectories D_k = {τ_i} by running policy π_k = π(θ_k) in the environment.
- 5: Compute rewards-to-go \hat{R}_t .
- Compute advantage estimates, Â_t (using any method of advantage estimation) based on the current value function V_{φ_t}.
- 7: Estimate policy gradient as

$$\hat{g}_k = \frac{1}{|\mathcal{D}_k|} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^T \nabla_{\theta} \log \pi_{\theta}(a_t | s_t)|_{\theta_k} \hat{A}_t.$$

8: Use the conjugate gradient algorithm to compute

$$\hat{x}_k \approx \hat{H}_k^{-1} \hat{g}_k$$
,

where \hat{H}_k is the Hessian of the sample average KL-divergence.

Update the policy by backtracking line search with

$$\theta_{k+1} = \theta_k + \alpha^j \sqrt{\frac{2\delta}{\hat{x}_k^T \hat{H}_k \hat{x}_k}} \hat{x}_k,$$

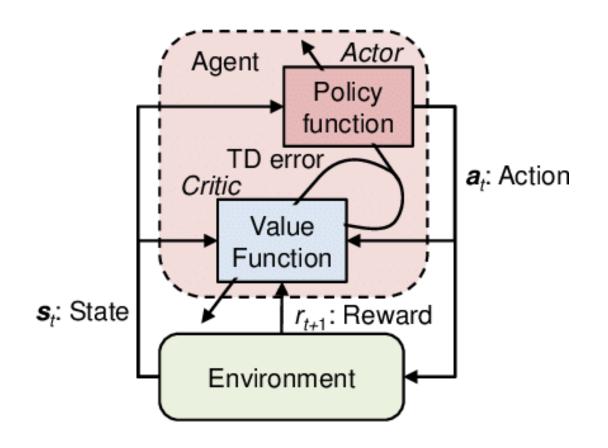
where $j \in \{0, 1, 2, ...K\}$ is the smallest value which improves the sample loss and satisfies the sample KL-divergence constraint.

0: Fit value function by regression on mean-squared error:

$$\phi_{k+1} = \arg \min_{\phi} \frac{1}{|\mathcal{D}_k|T} \sum_{\tau \in \mathcal{D}_k} \sum_{t=0}^{T} \left(V_{\phi}(s_t) - \hat{R}_t \right)^2,$$

typically via some gradient descent algorithm.

11: end for



Actor-Critic Method (A2C)

Problem: we don't know Q and V. Can we learn them?

Yes, using Q-learning! We can combine Policy Gradients and Q-learning by training both an **actor** (the policy) and a **critic** (the Q-function).

- The actor decides which action to take, and the critic tells the actor how good its action was and how it should adjust
- Also alleviates the task of the critic as it only has to learn the values of (state, action) pairs generated by the policy
- Can also incorporate Q-learning tricks e.g. experience replay
- Remark: we can define by the advantage function how much an action was better than expected $A^\pi(s,a) = Q^\pi(s,a) V^\pi(s)$

DDPG

- Deep Deterministic Policy Gradient (DDPG) is an algorithm which concurrently learns a Q-function and a policy. It uses off-policy data and the Bellman equation to learn the Q-function, and uses the Q-function to learn the policy
 - DDPG is an off-policy algorithm.
 - DDPG can only be used for environments with continuous action spaces.
 - DDPG can be thought of as being deep Q-learning for continuous action spaces.

• To make DDPG policies explore better, we add noise to their actions at training time. The authors of the original DDPG paper recommended time-correlated OU noise, but more recent results suggest that uncorrelated, mean-zero Gaussian noise works perfectly well.

Ornstein-Uhlenbeck process

RL - DDPG

Algorithm 1 Deep Deterministic Policy Gradient

- 1: Input: initial policy parameters θ , Q-function parameters ϕ , empty replay buffer \mathcal{D}
- 2: Set target parameters equal to main parameters $\theta_{\text{targ}} \leftarrow \theta$, $\phi_{\text{targ}} \leftarrow \phi$
- 3: repeat
- Observe state s and select action a = clip(μ_θ(s) + ε, a_{Low}, a_{High}), where ε ~ N
- Execute a in the environment
- 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}
- 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

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

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

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

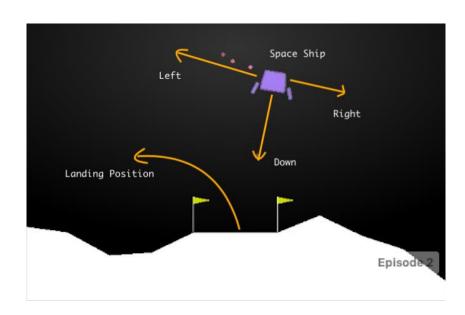
15: Update target networks with

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

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

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

Example



 Fuel is infinite, so an agent can learn to fly and then land on its first attempt.

Observation

- Landing pad is always at coordinates (0,0).
- Coordinates are the first two numbers in state vector.

Reward

- From the top of the screen to landing pad and zero speed is about 100..140 points.
- If lander moves away from landing pad it loses reward back.
- Episode finishes if the lander crashes or comes to rest, receiving additional -100 or +100 points.
- Each leg ground contact is +10.
- Firing main engine is -0.3 points each frame.
- Solved is 200 points.

Actions

Action is two real values vector from -1 to +1. First controls main engine, -1..0 off, 0..+1 throttle from 50% to 100% power. Engine can't work with less than 50% power. Second value -1.0..-0.5 fire left engine, +0.5..+1.0 fire right engine, -0.5..0.5 off.

Thanks

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