

Topics in Reinforcement Learning:  
Lessons from AlphaZero for  
(Sub)Optimal Control and Discrete Optimization

Arizona State University  
Course CSE 691, Spring 2022

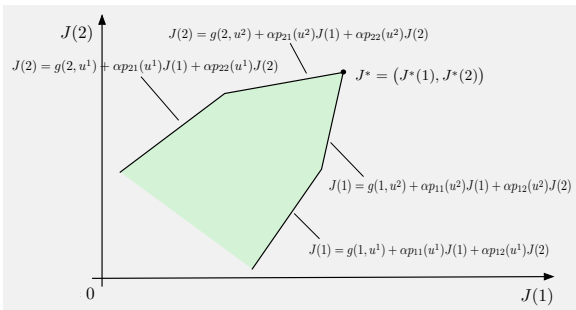
Links to Class Notes, Videolectures, and Slides at  
<http://web.mit.edu/dimitrib/www/RLbook.html>

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Lecture 11  
Approximate Linear Programming;  
Policy Gradient and Random Search Methods

- 1 Linear Programming: Another Approach to Approximation in Value Space
- 2 Approximation in Policy Space: Motivation
- 3 Training of Policies by Cost Optimization - Random Search
- 4 Training of Policies by Cost Optimization - Policy Gradient Methods
- 5 Implementation Issues of Policy Gradient Methods

# Exact Solution of Discounted DP by Linear Programming



**Key idea:**  $J^*$  is the “largest”  $J$  that satisfies the constraint

$$J(i) \leq \sum_{j=1}^n p_{ij}(u) (g(i, u, j) + \alpha J(j)), \quad \text{for all } i = 1, \dots, n \text{ and } u \in U(i),$$

so that  $J^* = (J^*(1), \dots, J^*(n))$  maximizes  $\sum_{i=1}^n J(i)$  subject to the above constraint.

**Proof:** Generate sequence  $\{J_k\}$  with VI, starting from any  $J = J_0$  satisfying the constraint, which implies that  $J_0 \leq J_1$ . Since  $J_k = T^k J_0$  and  $T$  is monotone, we have  $J = J_0 \leq J_k \leq J_{k+1} \rightarrow J^*$ . So any  $J$  satisfying the constraint also satisfies  $J \leq J^*$ .

Difficulty of the exact LP algorithm for large problems

**Too many variables** ( $n$ ) and **too many constraints** (the # of state-control pairs).

Introduce a linear feature-based architecture  $J^*(i) \approx \tilde{J}(i, r) = \sum_{\ell=1}^m r_{\ell} \phi_{\ell}(i)$

- **Replace  $J(i)$  with  $\tilde{J}(i, r)$**  to reduce the number of variables.
- **Introduce constraint sampling** to reduce the number of constraints.
- **Maximize  $\sum_{i \in \tilde{I}} \tilde{J}(i, r)$**  subject to

$$\tilde{J}(i, r) \leq \sum_{j=1}^n p_{ij}(u) (g(i, u, j) + \alpha \tilde{J}(j, r)), \quad i \in \tilde{I}, u \in \tilde{U}(i)$$

This is a linear program.

- $\tilde{I}$  is a set of “representative states”,  $\tilde{U}(i)$  is a set of “representative controls”.
- **Sampling with some known suboptimal policies is typically used to select a subset of the constraints to enforce**; progressively enrich the subset as necessary.
- The approach has not been used widely, but has been successful on substantive test problems (see Van Roy and De Farias’ works, among others).
- **Capitalizes on the reliability of large-scale LP software.**

# General Framework for Approximation in Policy Space

- **Parametrize stationary policies with a parameter vector  $r$** ; denote them by  $\tilde{\mu}(r)$ , with components  $\tilde{\mu}(i, r)$ ,  $i = 1, \dots, n$ . **Each  $r$  defines a policy.**
- The parametrization may be problem-specific, or feature-based, or may involve a neural network.
- The idea is to **optimize some measure of performance with respect to  $r$ .**

**An example of problem-specific/natural parametrization:** Supply chains, inventory control



- Retail center places orders to the production center, depending on current stock; there may be orders in transit; demand and delays can be stochastic.
- State is (current stock, orders in transit, ++). Can be formulated by DP but can be very difficult to solve exactly.
- Intuitively, a near-optimal policy is of the form: **When the retail inventory goes below level  $r_1$ , order an amount  $r_2$ . Optimize over the parameter vector  $r = (r_1, r_2)$ .**
- Extensions to a network of production/retail centers, multiple products, etc.

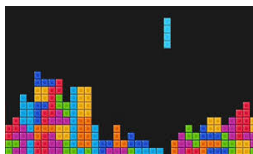
## Another Example: Policy Parametrization Through Value Parametrization

### Indirect parametrization of policies through cost features

- Suppose  $\tilde{J}(i, r)$  is a cost function parametric approximation.
- $\tilde{J}$  may be a linear feature-based architecture that is natural for the given problem.
- Define

$$\tilde{\mu}(i, r) \in \arg \min_{u \in U(i)} \sum_{j=1}^n p_{ij}(u) (g(i, u, j) + \tilde{J}(j, r))$$

- **This is useful when we know a good parametrization in value space**, but we want to use a method that works well in policy space, and results in an easily implementable policy.



**Tetris example:** There are good linear parametrizations through features. **Great success has been achieved by indirect approximation in policy space.**

Think about at least six contexts where approximation in policy space is either essential or is helpful

- Problems with natural policy parametrizations (like the supply chain problem)
- Problems with natural value parametrizations (like the tetris problem), where a good policy training method works well.
- Approximation in policy space on top of approximation in value space.
- Learning from a software or human expert.
- Unconventional information structures (limited memory, etc) - Conventional DP breaks down.
- Multiagent systems with local information (not shared with other agents).

- Compute approximate cost-to-go function  $\tilde{J}$  using an approximation in value space scheme.
- This defines the corresponding suboptimal policy  $\hat{\mu}$  through one-step lookahead,

$$\hat{\mu}(i, r) \in \arg \min_{u \in U(i)} \sum_{j=1}^n p_{ij}(u) (g(i, u, j) + \tilde{J}(j, r))$$

or a multistep lookahead version.

- Approximate  $\hat{\mu}$  using a training set consisting of a large number  $q$  of sample pairs  $(i^s, u^s)$ ,  $s = 1, \dots, q$ , where  $u^s = \hat{\mu}(i^s)$ .
- In particular, introduce a parametric family of policies  $\tilde{\mu}(i, r)$ . Then obtain  $r$  by

$$\min_r \sum_{s=1}^q \|u^s - \tilde{\mu}(i^s, r)\|^2.$$



- Suppose **we have a software or human expert** that can choose a “good” or “near-optimal” control  $u^s$  at any state  $i^s$ .
- We form a sample set of representative state-control pairs  $(i^s, u^s)$ ,  $s = 1, \dots, q$ .
- We introduce a parametric family of policies  $\tilde{\mu}(i, r)$ . Then obtain  $r$  by

$$\min_r \sum_{s=1}^q \|u^s - \tilde{\mu}(i^s, r)\|^2.$$

- This approach is known as **expert supervised training**.
- It has been used (in various forms) in backgammon and in chess.
- It can be used, among others, for initialization of other methods.

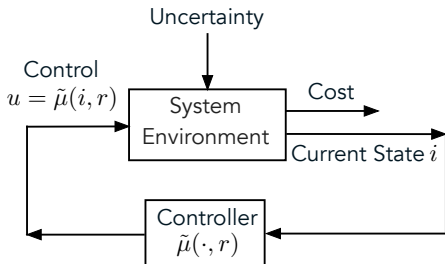
# Unconventional Information Structures

- Approximation in value space is based on a DP formulation, so the controller has access to the exact state (or a belief state in case of partial state information).
- In some contexts this may not be true. **There is a DP-like structure, but no full state or belief state is available.**
- **Example 1:** The controller “forgets” information, e.g., “limited memory”.
- **Example 2:** Some control components may be chosen on the basis of different information than others.

## Example: Multiagent systems with local agent information

- Suppose decision making and information gathering is distributed among multiple autonomous agents.
- **Each agent’s action depends only on his/her local information.**
- Agents may be receiving delayed information from other agents.
- Then **conventional DP and much of the approximation in value space methodology breaks down.**
- Approximation in policy space is still applicable.

# Optimization/Training Framework



## Training by Cost Optimization

- Each  $r$  defines a stationary policy  $\tilde{\mu}(r)$ , with components  $\tilde{\mu}(i, r)$ ,  $i = 1, \dots, n$ .
- Determine  $r$  through the minimization

$$\min_r J_{\tilde{\mu}(r)}(i_0)$$

where  $J_{\tilde{\mu}(r)}(i_0)$  is the cost of the policy  $\tilde{\mu}(r)$  starting from initial state  $i_0$ .

- More generally, determine  $r$  through the minimization

$$\min_r E\{J_{\tilde{\mu}(r)}(i_0)\}$$

where the  $E\{\cdot\}$  is with respect to a suitable probability distribution of  $i_0$ .

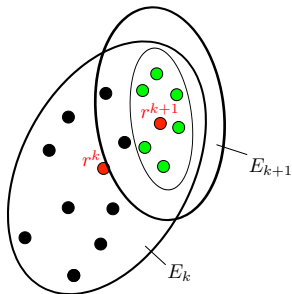
Random search methods apply to the general minimization  $\min_{r \in R} F(r)$

- They generate a parameter sequence  $\{r^k\}$  aiming for cost reduction.
- Given  $r^k$ , points are chosen in some random fashion in a neighborhood of  $r^k$ , and some new point  $r^{k+1}$  is chosen within this neighborhood.
- In theory they have good convergence properties. In practice they can be slow.
- They are not affected as much by local minima (as for example gradient-type methods).
- They don't require a differentiable cost function, and they apply to discrete as well as continuous minimization.
- There are many methods and variations thereof.

## Some examples

- Evolutionary programming.
- Tabu search.
- Simulated annealing.
- Cross entropy method.

## Cross-Entropy Method - A Sketch



- At the current iterate  $r^k$ , construct an ellipsoid  $E_k$  centered at  $r^k$ .
- Generate a number of random samples within  $E_k$ . “Accept” a subset of the samples that have “low” cost.
- Let  $r^{k+1}$  be the sample “mean” of the accepted samples.
- Construct a sample “covariance” matrix of the accepted samples, form the new ellipsoid  $E_{k+1}$  using this matrix, and continue.
- Limited convergence rate guarantees. Success depends on domain-specific insight and the skilled use of implementation heuristics.
- Simple and well-suited for parallel computation. **Resembles a “gradient method”.**

Consider the minimization of  $J_{\tilde{\mu}(r)}(i_0)$  over  $r$  by using the gradient method

$$r^{k+1} = r^k - \gamma^k \nabla J_{\tilde{\mu}(r^k)}(i_0)$$

assuming that  $J_{\tilde{\mu}(r)}(i_0)$  is differentiable with respect to  $r$ .

- The difficulty is that the gradient  $\nabla J_{\tilde{\mu}(r^k)}(i_0)$  may not be explicitly available.
- Then the gradient must be approximated by finite differences of cost function values  $J_{\tilde{\mu}(r^k)}(i_0)$ .
- **When the problem is deterministic the gradient method may work well.**
- **When the problem is stochastic**, the cost function values may be computable only through Monte Carlo simulation. **Very hard to get accurate gradients by differencing function values.**

Consider the generic optimization problem  $\min_{z \in Z} F(z)$

We take an unusual step: **Convert this problem to the stochastic optimization problem**

$$\min_{p \in \mathcal{P}_Z} E_p\{F(z)\}$$

where

- $z$  is viewed as a random variable.
- $\mathcal{P}_Z$  is the set of probability distributions over  $Z$ .
- $p$  denotes the generic distribution in  $\mathcal{P}_Z$ .
- $E_p\{\cdot\}$  denotes expected value with respect to  $p$ .

How does this relate to our infinite horizon DP problems?

- For this framework to apply to a stochastic DP context, **we must enlarge the set of policies to include randomized policies**, mapping a state  $i$  into a probability distribution over the set of controls  $U(i)$ .
- Note that in our DP problems, **optimization over randomized policies gives the same results as optimization over ordinary/nonrandomized policies**.
- In the DP context,  $z$  is the state-control trajectory:  $z = \{i_0, u_0, i_1, u_1, \dots\}$ .

## Parametrization of the probability distributions

- We restrict attention to a parametrized subset  $\tilde{\mathcal{P}}_Z \subset \mathcal{P}_Z$  of probability distributions  $p(z; r)$ , where  $r$  is a continuous parameter.
- In other words, we approximate the problem  $\min_{z \in Z} F(z)$  with **the restricted problem**

$$\min_r E_{p(z;r)}\{F(z)\}$$

- We use a gradient method for solving this problem:

$$r^{k+1} = r^k - \gamma^k \nabla \left( E_{p(z;r^k)}\{F(z)\} \right)$$

- Key fact: **There is a useful formula for the gradient**, which involves the gradient with respect to  $r$  of the natural logarithm  $\log(p(z; r^k))$ .



Assuming that  $p(z; r^k)$  is a discrete distribution, we have

$$\begin{aligned}\nabla\left(E_{p(z; r^k)}\{F(z)\}\right) &= \nabla\left(\sum_{z \in Z} p(z; r^k) F(z)\right) \\ &= \sum_{z \in Z} \nabla p(z; r^k) F(z) \\ &= \sum_{z \in Z} p(z; r^k) \frac{\nabla p(z; r^k)}{p(z; r^k)} F(z) \\ &= E_{p(z; r^k)}\left\{\nabla\left(\log(p(z; r^k))\right) F(z)\right\}\end{aligned}$$

Sample-Based Gradient Method for Parametric Approximation of  $\min_{z \in Z} F(z)$

- At  $r^k$  obtain a sample  $z^k$  according to the distribution  $p(z; r^k)$ .
- Compute the sample gradient  $\nabla\left(\log(p(z^k; r^k))\right) F(z^k)$ .
- Use it to iterate according to

$$r^{k+1} = r^k - \gamma^k \nabla\left(\log(p(z^k; r^k))\right) F(z^k)$$

- Denote by  $z$  the infinite horizon state-control trajectory:

$$z = \{i_0, u_0, i_1, u_1, \dots\}.$$

- We consider a **parametrization of randomized policies**  $p(u | i; r)$  with parameter  $r$ , i.e., the control at state  $i$  is generated according to a distribution  $p(u | i; r)$  over  $U(i)$ .
- Then for a given  $r$ , the state-control trajectory  $z$  is a random trajectory with probability distribution denoted  $p(z; r)$ .
- The cost corresponding to the trajectory  $z$  is

$$F(z) = \sum_{m=0}^{\infty} \alpha^m g(i_m, u_m, i_{m+1}),$$

and the problem is to minimize  $E_{p(z;r)}\{F(z)\}$ , over  $r$ .

- The gradient needed in the gradient iteration

$$r^{k+1} = r^k - \gamma^k \nabla \left( \log(p(z^k; r^k)) \right) F(z^k)$$

is given by

$$\nabla \left( \log(p(z^k; r^k)) \right) = \sum_{m=0}^{\infty} \log(p_{i_m i_{m+1}}(u_m)) + \sum_{m=0}^{\infty} \nabla \left( \log(p(u_m | i_m; r^k)) \right)$$

- It involves the cost function of the discounted problem, but not its gradient ... In fact the cost per stage  $g$  may be nondifferentiable!
- The problem solved is a randomized version of the original ... so if  $r^k \rightarrow \bar{r}$  and the distribution  $p(z, \bar{r})$  is not atomic, a solution has to be extracted from this distribution.

## Some of the implementation issues

- How to collect the trajectory samples  $z^k$  to strike a **balance between convenient implementation and exploration** of the search space.
- How to **reduce the large noise** in the cost calculation  $F(z^k)$ .
- Use of **baseline  $b$** , i.e., iterate according to

$$r^{k+1} = r^k - \gamma^k \nabla \left( \log(p(z^k; r^k)) \right) (F(z^k) - b)$$

instead of

$$r^{k+1} = r^k - \gamma^k \nabla \left( \log(p(z^k; r^k)) \right) F(z^k)$$

There is theoretical basis for this (see the next slide).

Introduce an equivalent “variational” problem (known since the 1960s)

- Subtract any known function  $V(x)$  from  $J^*(x)$ :

$$\hat{J}(x) = J^*(x) - V(x), \quad x = 1, \dots, n$$

- Replace the cost per stage  $g(x, u, y)$  with

$$\hat{g}(x, u, y) = g(x, u, y) + \alpha V(y) - V(x), \quad x = 1, \dots, n$$

- Then the original problem’s Bellman’s equation is written as another Bellman equation

$$\hat{J}(x) = \min_{u \in U(x)} \sum_{y=1}^n p_{xy}(u) (\hat{g}(x, u, y) + \alpha \hat{J}(y)), \quad x = 1, \dots, n$$

- $\hat{J}$  is the optimal cost of another problem:  $g(x, u, y)$  is replaced by  $\hat{g}(x, u, y)$
- The reformulated problem is equivalent as far as exact solution is concerned
- BUT  $\hat{J}$  may have more favorable “shape” for approximation, i.e., policy gradient and other methods may work better for the reformulated problem
- **Example:** If  $V \approx J^*$ , approximation methods can capture more easily small scale variations in  $J^*$  ... compare with the discussion on advantage updating (Lecture 8)

There is a generic difficulty with using a fixed policy on-line:

- It is **all-training no on-line play**. (This could be good but could be very bad.)
- **It does not adapt to changes in the problem's parameters**.
- So approximation in policy space may not work well in adaptive control contexts.
- Also it does not yield the benefit of on-line lookahead minimization/rollout.
- Approximation in value space, and rollout may work much better (e.g., in **AlphaZero**).

An alternative use of approximation in policy space methods (including policy gradient)

It can provide a base policy for use in (truncated) rollout or can be used in Monte Carlo Tree Search. This is what is done in AlphaZero.

## About the Next Lecture

We will cover aggregation, which trains off-line a cost function approximation.

We will use videolecture 12 from the 2021 ASU class.