

# A Series of Lectures on Approximate Dynamic Programming

## Lecture 2

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# APPROXIMATE DYNAMIC PROGRAMMING I

- 1 Approximation in Value Space - Limited Lookahead
- 2 Parametric Cost Approximation

## Recall our Problem Structure

Discrete-time system

$$x_{k+1} = f_k(x_k, u_k, w_k), \quad k = 0, 1, \dots, N-1$$

- $x_k$ : State
- $u_k$ : Control from a constraint set  $U_k(x_k)$
- $w_k$ : Disturbance; random parameter with distribution  $P(w_k | x_k, u_k)$

Optimization over Feedback Policies  $\pi = \{\mu_0, \mu_1, \dots, \mu_{N-1}\}$ , with  
 $u_k = \mu_k(x_k) \in U(x_k)$

Cost of a policy starting at initial state  $x_0$ :

$$J_\pi(x_0) = E \left\{ g_N(x_N) + \sum_{k=0}^{N-1} g_k(x_k, \mu_k(x_k), w_k) \right\}$$

Optimal cost function:

$$J^*(x_0) = \min_{\pi} J_{\pi}(x_0)$$

## Recall the Exact DP Algorithm

Computes for all  $k$  and states  $x_k$ :  $J_k(x_k)$ , the opt. cost of tail problem that starts at  $x_k$

Go backwards,  $k = N - 1, \dots, 0$ , using

$$J_N(x_N) = g_N(x_N)$$

$$J_k(x_k) = \min_{u_k \in U_k(x_k)} E\left\{g_k(x_k, u_k, w_k) + J_{k+1}(f_k(x_k, u_k, w_k))\right\}$$

### Notes:

- $J_0(x_0) = J^*(x_0)$ : Cost generated at the last step, is equal to the optimal cost
- Let  $\mu_k^*(x_k)$  minimize in the right side above for each  $x_k$  and  $k$ . Then the policy  $\pi^* = \{\mu_0^*, \dots, \mu_{N-1}^*\}$  is optimal

## Practical Difficulties of DP

### The curse of dimensionality (too many values of $x_k$ )

- In continuous-state problems:
  - ▶ Discretization needed
  - ▶ Exponential growth of the computation with the dimensions of the state and control spaces
- In naturally discrete/combinatorial problems: Quick explosion of the number of states as the search space increases
- Length of the horizon (what if it is infinite?)

### The curse of modeling; we may not know exactly $f_k$ and $P(x_k | x_k, u_k)$

- It is often hard to construct an accurate math model of the problem
- Sometimes a simulator of the system is easier to construct than a model

### The problem data may not be known well in advance

- A family of problems may be addressed. The data of the problem to be solved is given with little advance notice
- The problem data may change as the system is controlled – need on-line replanning and fast solution

# A MAJOR IDEA: Cost Approximation

One-Step Lookahead - Idea is to simplify the DP computation

- Replace  $J_{k+1}$  by an approximation  $\tilde{J}_{k+1}$
- Apply  $\bar{u}_k$  that attains the minimum in

$$\min_{u_k \in U_k(x_k)} E \left\{ g_k(x_k, u_k, w_k) + \tilde{J}_{k+1}(f_k(x_k, u_k, w_k)) \right\}$$

$\ell$ -Step Lookahead

- At state  $x_k$  solve the  $\ell$ -step DP problem starting at  $x_k$  and using terminal cost  $\tilde{J}_{k+\ell}$
- If  $\bar{u}_k, \bar{u}_{k+1}, \dots, \bar{u}_{k+\ell-1}$  is an optimal policy for the  $\ell$ -step problem, apply the first control  $\bar{u}_k$

Other Names Used

Rolling or receding horizon control

Let's focus on the one-step lookahead computation at stage  $k$

$$\min_{u_k \in U_k(x_k)} E\left\{ g_k(x_k, u_k, w_k) + \tilde{J}_{k+1}(f_k(x_k, u_k, w_k)) \right\}$$

## Issues

- A key issue: How do we choose the approximate cost functions  $\tilde{J}_k$ ?
- Another issue: How do we deal with the minimization and the computation of  $E\{\cdot\}$

A variety of approximation approaches (and combinations thereof):

- **Parametric cost-to-go approximation:** Use as  $\tilde{J}_k$  a parametric function  $\tilde{J}_k(x_k, r_k)$  (e.g., a neural network), whose parameter  $r_k$  is "tuned" by some scheme
- **Rollout:** Use as  $\tilde{J}_k$  the cost of some suboptimal policy, which is calculated either analytically or by simulation
- **Problem approximation:** Use  $\tilde{J}_k$  derived from a related but simpler problem

At State  $x_k$

DP minimization

$$\min_{u_k, \mu_{k+1}, \dots, \mu_{k+\ell-1}} E \left\{ g_k(x_k, u_k, w_k) + \sum_{m=k+1}^{k+\ell-1} g_k(x_m, \mu_m(x_m), w_m) + \tilde{J}_{k+\ell}(x_{k+\ell}) \right\}$$

First  $\ell$  Steps

“Future”

Lookahead Minimization

Cost-to-go Approximation

**At State  $x_k$**   
**DP minimization**

$$\min_{u_k, \mu_{k+1}, \dots, \mu_{k+\ell-1}} E \left\{ g_k(x_k, u_k, w_k) + \sum_{m=k+1}^{k+\ell-1} g_k(x_m, \mu_m(x_m), w_m) + \tilde{J}_{k+\ell}(x_{k+\ell}) \right\}$$

First  $\ell$  Steps      “Future”

↑

Computation of  $\tilde{J}_{k+\ell}$ :

- Simple choices
- Parametric approximation
- Rollout
- Tail problem approximation

**At State  $x_k$**

**DP minimization**

$$\min_{u_k, \mu_{k+1}, \dots, \mu_{k+\ell-1}} E \left\{ g_k(x_k, u_k, w_k) + \sum_{m=k+1}^{k+\ell-1} g_k(x_m, \mu_m(x_m), w_m) + \tilde{J}_{k+\ell}(x_{k+\ell}) \right\}$$

**Approximations:**

Replace  $E\{\cdot\}$  with nominal values  
(certainty equivalent control)

Limited simulation  
(Monte Carlo tree search)

**Computation of  $\tilde{J}_{k+\ell}$ :**

- Simple choices
- Parametric approximation
- Rollout
- Tail problem approximation

## Long lookahead $\ell$ and simple choice of $\tilde{J}_{k+\ell}$

- Some examples

$$\tilde{J}_{k+\ell}(x) \equiv 0 \quad (\text{or a constant})$$

$$\tilde{J}_{k+\ell}(x) = g_N(x)$$

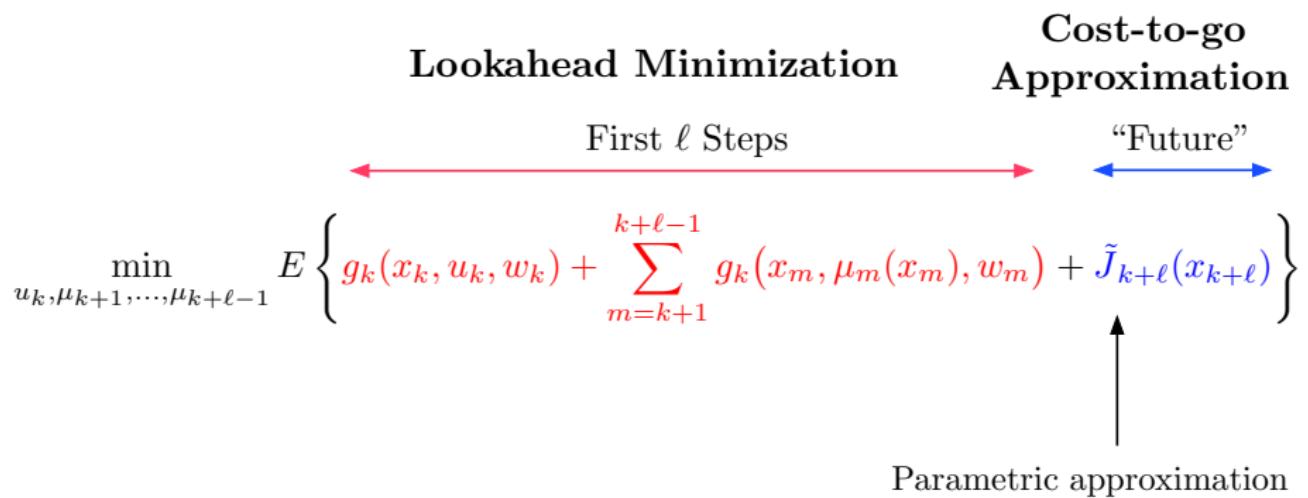
For problems with a "goal state" use a simple penalty  $\tilde{J}_{k+\ell}$

$$\tilde{J}_{k+\ell}(x) = \begin{cases} 0 & \text{if } x \text{ is a goal state} \\ >> 1 & \text{if } x \text{ is not a goal state} \end{cases}$$

- Long lookahead  $\implies$  A lot of DP computation
- Often must be done off-line

## Short lookahead $\ell$ and sophisticated choice $\tilde{J}_{k+\ell} \approx J_{k+\ell}$

- The lookahead cost function approximates (to within a constant) the optimal cost-to-go produced by exact DP
- We will next describe a variety of off-line and on-line approximation approaches



$$J_k(x_k) \approx \tilde{J}_k(x_k, r_k)$$

with

$r_k = (r_{1,k}, \dots, r_{m,k})$  a vector of "tunable" scalar weights

- We use  $\tilde{J}_k$  in place of  $J_k$  (the optimal cost-to-go function) in a one-step or multistep lookahead scheme
- $\tilde{J}_k(x_k, r_k)$  is called an **approximation architecture**
- **Role of  $r_k$ :** By adjusting  $r_k$  we can change the "shape" of  $\tilde{J}_k$  so that it is "close" to the optimal  $J_k$  (at least within a constant)

## Two key Issues

- The choice of the parametric class  $\tilde{J}_k(x_k, r_k)$ ; there is a large variety
- The method for tuning/adjusting the weights ("training" the architecture)

## Feature extraction

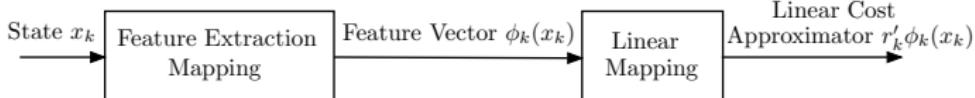
- A process that maps the state  $x_k$  into a vector  $\phi_k(x_k) = (\phi_{1,k}(x_k), \dots, \phi_{m,k}(x_k))$ , called the **feature vector** associated with  $x_k$
- A feature-based cost approximator has the form

$$\tilde{J}_k(x_k, r_k) = \hat{J}_k(\phi_k(x_k), r_k)$$

where  $r_k$  is a parameter vector and  $\hat{J}_k$  is some function, linear or nonlinear in  $r_k$

- With a well-chosen feature vector  $\phi_k(x_k)$ , a good approximation to the cost-to-go is often provided by **linearly** weighting the features, i.e.,

$$\tilde{J}_k(x_k, r_k) = \hat{J}_k(\phi_k(x_k), r_k) = \sum_{i=1}^m r_{i,k} \phi_{i,k}(x_k) = r'_k \phi_k(x_k)$$

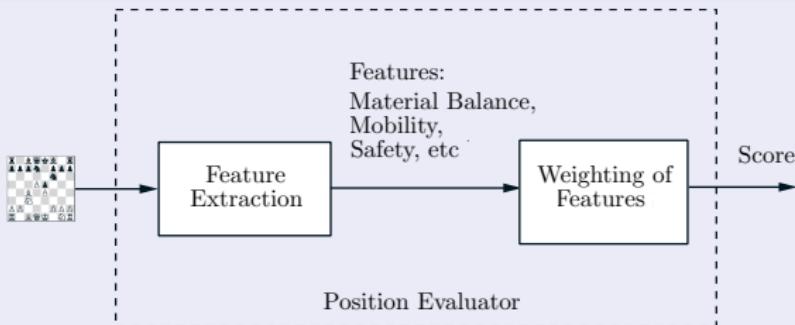


This can be viewed as **subspace approximation**; view the features the  $\phi_{i,k}(x_k)$  as basis functions

# Feature Selection: A Major Issue

- Any generic basis functions, such as classes of polynomials, wavelets, radial basis functions, etc, can serve as features
- In some cases, problem-specific features can be “hand-crafted”

## Computer chess example



- Think of **state**: board position; **control**: move choice
- Use a **feature-based position evaluator assigning a score to each position**
- Most chess programs use a linear architecture with “manual” choice of weights
- Some computer programs choose the weights by a least squares fit using lots of grandmaster play examples

# An Example of Architecture Training: Sequential DP Approximation

A common way to train architectures  $\tilde{J}_k(x_k, r_k)$  in the context of DP

- We start with  $\tilde{J}_N = g_N$  and sequentially train going backwards, until  $k = 1$
- Given a cost-to-go approximation  $\tilde{J}_{k+1}$ , we use one-step lookahead to construct a large number of state-cost pairs  $(x_k^s, \beta_k^s)$ ,  $s = 1, \dots, q$ , where

$$\beta_k^s = \min_{u \in U_k(x_k^s)} E \left\{ g(x_k^s, u, w_k) + \tilde{J}_{k+1} \left( f_k(x_k^s, u, w_k), r_{k+1} \right) \right\}, \quad s = 1, \dots, q$$

- We "train" an architecture  $\tilde{J}_k$  on the training set  $(x_k^s, \beta_k^s)$ ,  $s = 1, \dots, q$

## Training by least squares/regression

- We minimize over  $r_k$

$$\sum_{s=1}^q (\tilde{J}_k(x_k^s, r_k) - \beta_k^s)^2 + \gamma \|r_k - \bar{r}\|^2$$

where  $\bar{r}$  is an initial guess for  $r_k$  and  $\gamma > 0$  is a regularization parameter

- Incremental gradient methods are typically used for this. They take advantage of the large sum structure of the cost function
- For a linear architecture the training problem is a linear least squares problem

Neural nets can be used in a sequential DP approximation scheme: Train the stage  $k$  neural net (i.e., compute  $\tilde{J}_k$ ) using a training set generated with the stage  $k + 1$  neural net (which defines  $\tilde{J}_{k+1}$ )

Focus at the typical stage  $k$  and drop the index  $k$  for convenience

- Neural nets are approximation architectures of the form

$$\tilde{J}(x, v, r) = \sum_{i=1}^m r_i \phi_i(x, v) = r' \phi(x, v)$$

involving two parameter vectors  $r$  and  $v$  with different roles

- View  $\phi(x, v)$  as a feature vector; view  $r$  as a vector of linear weighting parameters for  $\phi(x, v)$
- The training is done by least squares/regression
- By training  $v$  jointly with  $r$ , we obtain automatically generated features!