MIT OpenCourseWare <http://ocw.mit.edu>

6.231 Dynamic Programming and Stochastic Control Fall 2008

For information about citing these materials or our Terms of Use, visit: [http://ocw.mit.edu/terms.](http://ocw.mit.edu/terms)

## 6.231 DYNAMIC PROGRAMMING

### LECTURE 12

## LECTURE OUTLINE

- More on rollout algorithms Stochastic problems
- Simulation-based methods for rollout
- Approximations of rollout algorithms
- Rolling horizon approximations
- Discretization of continuous time
- Discretization of continuous space
- Other suboptimal approaches

#### ROLLOUTALGORITHMS -STOCHASTIC PROBLEM

**Rollout policy:** At each k and state  $x_k$ , use the control  $\overline{\mu}_k(x_k)$  that

$$
\min_{u_k \in U_k(x_k)} Q_k(x_k, u_k),
$$

where

$$
Q_k(x_k, u_k) = E\big\{g_k(x_k, u_k, w_k) + H_{k+1}\big(f_k(x_k, u_k, w_k)\big)\big\}
$$

and  $H_{k+1}(x_{k+1})$  is the cost-to-go of the heuristic.

 $Q_k(x_k, u_k)$  is called the *Q-factor* of  $(x_k, u_k)$ , and for a stochastic problem, its computation may involve Monte Carlo simulation.

• Potential difficulty: To minimize over  $u_k$  the  $Q$ factor, we must form Q-factor differences  $Q_k(x_k, u)$ −  $Q_k(x_k, \overline{u})$ . This differencing often amplifies the simulation error in the calculation of the Q-factors.

Potential remedy: Compare any two controls u and  $\overline{u}$  by simulating the *Q-factor differences*  $Q_k(x_k, u) - Q_k(x_k, \overline{u})$  directly. This may effect variance reduction of the simulation-induced error.

# Q-FACTOR APPROXIMATION

Here, instead of simulating the  $Q$ -factors, we approximate the costs-to-go  $H_{k+1}(x_{k+1})$ .

Certainty equivalence approach: Given  $x_k$ , fix future disturbances at "typical" values  $\overline{w}_{k+1}, \ldots, \overline{w}_{N-1}$ and approximate the Q-factors with

$$
\tilde{Q}_k(x_k, u_k) = E\big\{g_k(x_k, u_k, w_k) + \tilde{H}_{k+1}\big(f_k(x_k, u_k, w_k)\big)\big\}
$$

where  $\tilde{H}_{k+1}(f_k(x_k, u_k, w_k))$  is the cost of the heuristic with the disturbances fixed at the typical values.

• This is an approximation of  $H_{k+1}(f_k(x_k, u_k, w_k))$ by using a "single sample simulation."

• Variant of the certainty equivalence approach: Approximate  $H_{k+1}(f_k(x_k, u_k, w_k))$  by simulation using a small number of "representative samples" (scenarios).

• Alternative: Calculate (exact or approximate) values for the cost-to-go of the base policy at a limited set of state-time pairs, and then approximate  $H_{k+1}$  using an approximation architecture and a "training algorithm" or "least-squares fit."

## ROLLING HORIZON APPROACH

This is an *l*-step lookahead policy where the cost-to-go approximation is just 0.

• Alternatively, the cost-to-go approximation is the terminal cost function  $q_N$ .

• A short rolling horizon saves computation.

• "Paradox": It is not true that a longer rolling horizon always improves performance.

Example: At the initial state, there are two controls available (1 and 2). At every other state, there is only one control.



### ROLLING HORIZON COMBINED WITH ROLLOUT

We can use a rolling horizon approximation in calculating the cost-to-go of the base heuristic.

• Because the heuristic is suboptimal, the rationale for a long rolling horizon becomes weaker.

**Example:**  $N$ -stage stopping problem where the stopping cost is 0, the continuation cost is either  $-\epsilon$  or 1, where  $0 < \epsilon \ll 1$ , and the first state with continuation cost equal to 1 is state  $m$ . Then the optimal policy is to stop at state  $m$ , and the optimal cost is  $-m\epsilon$ .



Consider the heuristic that continues at every state, and the rollout policy that is based on this heuristic, with a rolling horizon of  $l \leq m$  steps.

It will continue up to the first  $m - l + 1$  stages, thus compiling a cost of  $-(m-l+1)\epsilon$ . The rollout performance improves as l becomes shorter!

Limited vision may work to our advantage!

# DISCRETIZATION

If the state space and/or control space is continuous/infinite, it must be replaced by a finite discretization.

Need for consistency, i.e., as the discretization becomes finer, the cost-to-go functions of the discretized problem converge to those of the continuous problem.

• Pitfalls with discretizing continuous time.

The control constraint set changes a lot as we pass to the discrete-time approximation.

• Continuous-Time Shortest Path Pitfall:

$$
\dot{x}_1(t) = u_1(t), \qquad \dot{x}_2(t) = u_2(t),
$$

with control constraint  $u_i(t) \in \{-1,1\}$  and cost  $\int_0^T g(x(t)) dt$ . Compare with naive discretization  $x_1(t+\Delta t) = x_1(t) + \Delta t u_1(t), x_2(t+\Delta t) = x_2(t) + \Delta t u_2(t)$ with  $u_i(t) \in \{-1, 1\}.$ 

• "Convexification effect" of continuous time.

#### SPACE DISCRETIZATION I

Given a discrete-time system with state space S, consider a finite subset  $\overline{S}$ ; for example  $\overline{S}$  could be a finite grid within a continuous state space S.

Difficulty:  $f(x, u, w) \notin \overline{S}$  for  $x \in \overline{S}$ .

We define an approximation to the original problem, with state space  $\overline{S}$ , as follows:

• Express each  $x \in S$  as a convex combination of states in  $\overline{S}$ , i.e.,

$$
x = \sum_{x_i \in \overline{S}} \gamma_i(x) x_i \quad \text{where } \gamma_i(x) \ge 0, \sum_i \gamma_i(x) = 1
$$

Define a "reduced" dynamic system with state space  $\overline{S}$ , whereby from each  $x_i \in \overline{S}$  we move to  $\overline{x} = f(x_i, u, w)$  according to the system equation of the original problem, and then move to  $x_j \in \overline{S}$ with probabilities  $\gamma_i(\overline{x})$ .

• Define similarly the corresponding cost per stage of the transitions of the reduced system.

## SPACE DISCRETIZATION II

• Let  $\overline{J}_k(x_i)$  be the optimal cost-to-go of the "reduced" problem from each state  $x_i \in \overline{S}$  and time k onward.

Approximate the optimal cost-to-go of any  $x \in$ S for the original problem by

$$
\tilde{J}_k(x) = \sum_{x_i \in \overline{S}} \gamma_i(x) \overline{J}_k(x_i),
$$

and use one-step-lookahead based on  $\tilde{J}_k$ .

The choice of coefficients  $\gamma_i(x)$  is in principle arbitrary, but should aim at consistency, i.e., as the number of states in  $\overline{S}$  increases,  $\widetilde{J}_k(x)$  should converge to the optimal cost-to-go of the original problem.

• Interesting observation: While the original problem may be deterministic, the reduced problem is always stochastic.

• Generalization: The set  $\overline{S}$  may be any finite set (not a subset of S) as long as the coefficients  $\gamma_i(x)$ admit a meaningful interpretation that quantifies the degree of association of x with  $x_i$ .

### OTHER SUBOPTIMAL CONTROL APPROACHES

• Minimize the DP equation error: Approximate the optimal cost-to-go functions  $J_k(x_k)$  with functions  $\tilde{J}_k(x_k, r_k)$ , where  $r_k$  is a vector of unknown parameters, chosen to minimize some form of error in the DP equations.

Direct approximation of control policies: For a subset of states  $x^i$ ,  $i = 1, \ldots, m$ , find

$$
\hat{\mu}_k(x^i) = \arg \min_{u_k \in U_k(x^i)} E\{g(x^i, u_k, w_k) + \tilde{J}_{k+1}(f_k(x^i, u_k, w_k), r_{k+1})\}.
$$

Then find  $\tilde{\mu}_k(x_k, s_k)$ , where  $s_k$  is a vector of parameters obtained by solving the problem

$$
\min_{s} \sum_{i=1}^{m} \|\hat{\mu}_k(x^i) - \tilde{\mu}_k(x^i, s)\|^2.
$$

• Approximation in policy space: Do not bother with cost-to-go approximations. Parametrize the policies as  $\tilde{\mu}_k(x_k, s_k)$ , and minimize the cost function of the problem over the parameters  $s_k$ .