## **Approximation in Value Space**

Let's begin by writing the DP algorithm:

$$J_N(x_n) = g_N(x_N)$$
 
$$J_k(x_k) = \min_{u_k \in U_k(x_k)} \left\{ \mathbb{E}_{w_k} \left[ g_k(x_k, u_k, w_k) + J_{k+1}(f_k(x_k, u_k, w_k)) \right] \right\}, \forall k \in \{0, ..., N-1\}$$

• Where we can compute a sub-optimal admissible policy by using an approximate cost-to-go function  $\tilde{J}_{k+1}(x_{k+1})$ :

$$\tilde{\mu}_k(x_k) \in \arg\min_{u_k \in U_k(x_k)} \left\{ \mathbb{E}_{w_k} \left[ g_k(x_k, u_k, w_k) + \tilde{J}_{k+1}(f_k(x_k, u_k, w_k)) \right] \right\}, \forall k \in \{0, ..., N-1\}$$

•  $\tilde{J}_{k+1}(x_{k+1})$  is also called an **approximate value function**.

### Multistep Lookahead

• As we saw last time, we can let the approximate function  $\tilde{J}_{k+1}(x_{k+1})$  be itself a l-step version of the DP problem:

$$\tilde{J}_{k+1}(x_{k+1}) = \min_{(\mu_{k+1}, \dots, \mu_{k+l-1})} \mathbb{E}_{w_{k+1}, \dots, w_{k+l-1}} \left[ \tilde{J}_{k+l}(x_{k+l}) + \sum_{i=k+1}^{\kappa_{k+l-1}} g_i(x_i, \mu_i(x_i), w_i) \right]$$

- Where  $\tilde{J}_{k+l}(x_{k+l})$  is yet another (albeit simpler) approximate value function.
- The multi-step Lookahead can be summarized as follows:

$$\min_{u_k,\mu_{k+1},...,\mu_{k+l-1}} \mathbb{E}_{w_{k+1},...,w_{k+l-1}} \left[ g_k(x_k,u_k,w_k) + \sum_{i=k+1}^{\text{First $l$-steps}} g_i(x_i,\mu_i(x_i),w_i) + \tilde{J}_{k+l}(x_{k+l}) \right]$$
 DP Minimization 
$$\text{Lookahead minimization} \qquad \text{Cost-to-go Approximation}$$

# **Multistep Lookahead**

• Ideally the lookahead length is "large-enough" so that we may use very simple approximate functions at the end of the lookahead period. For example:

$$\tilde{J}_{k+l}(x_{k+l}) \equiv 0$$
, or  $\tilde{J}_{k+l}(x_{k+l}) = g_N(x_{k+l})$ 

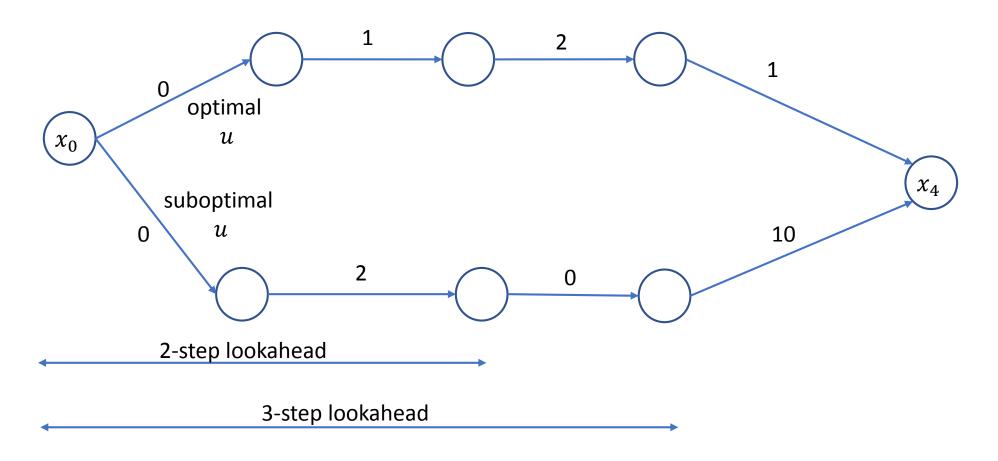
• In the multistep lookahead we compute a single control  $u_k$  that get's applied at stage k, and a sequence of functions  $(\mu_{k+1}(x_{k+1}), \dots, \mu_{k+l-1}(x_{k+l-1}))$  which gets **discarded.** 

• After applying  $u_k$  and obtaining the next state  $x_{k+1}$ , the whole system "rolls forward" and we have to re-solve the multistep-lookahead problem.

• This approach is called the **Rolling Horizon** approach, which is quite intuitive.

### **Example: Pitfall of Multistep Lookahead**

Consider this simple 4-stage Shortest-path problem:



So, a larger lookahead window may decrease performance.

### **Example: Partially Deterministic Lookahead**

• The multistep lookahead is very flexible. Consider again the I-step problem:

$$\tilde{\mu}_{k}(x_{k}) \in \arg\min_{u_{k} \in U_{k}(x_{k})} \left\{ \mathbb{E}_{w_{k}} \left[ g_{k}(x_{k}, u_{k}, w_{k}) + \tilde{J}_{k+1}(f_{k}(x_{k}, u_{k}, w_{k})) \right] \right\}, \forall k \in \{0, ..., N-1\}$$

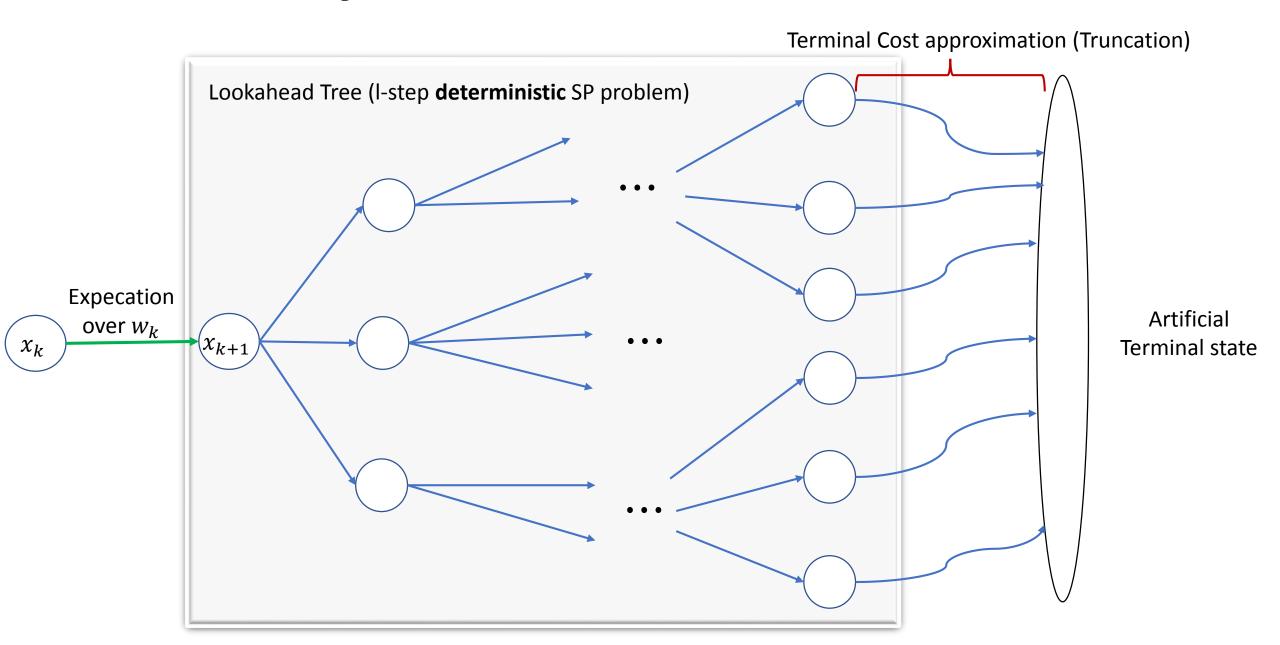
$$\tilde{J}_{k+1}(x_{k+1}) = \min_{(\mu_{k+1}, ..., \mu_{k+l-1})} \mathbb{E}_{w_{k+1}, ..., w_{k+l-1}} \left[ \tilde{J}_{k+l}(x_{k+l}) + \sum_{i=k+1}^{k+l-1} g_{i}(x_{i}, \mu_{i}(x_{i}), w_{i}) \right]$$

• Suppose that we apply the *Certainty Equivalence* principle where we replace the disturbances  $(w_{k+1}, ..., w_{k+l})$  by their average ("typical") values  $(\overline{w}_{k+1}, ..., \overline{w}_{k+l})$ :

$$\bar{w}_{k+1} = \mathbb{E}[w_{k+1}|x_{k+1}, u_{k+1}]$$

• Then the approximate value function  $\tilde{J}_{k+1}(x_{k+1})$  can be computed by a Deterministic Shortest Path Algorithm.

# **Partially Deterministic Lookahead**



#### **Scenario-based Lookahead**

 Note that the Certainty Equivalence approach can be generalized by using multiple scenarios in computing the expectation in:

$$\tilde{J}_{k+1}(x_{k+1}) = \min_{(\mu_{k+1}, \dots, \mu_{k+l-1})} \mathbb{E}_{w_{k+1}, \dots, w_{k+l-1}} \left[ \tilde{J}_{k+l}(x_{k+l}) + \sum_{i=k+1}^{k+l-1} g_i(x_i, \mu_i(x_i), w_i) \right]$$

• Suppose we have in our disposal a simulator that is able to generate a *scenario* sequence given the state  $x_{k+1}$ :

$$w^{s}(x_{k+1}) = (w_{k+1}^{s}, ..., w_{k+l-1}^{s}), s \in \{1, ..., S\}$$

• And let  $C^s(x_{k+1})$  be the solution of the **deterministic lookahead problem** under scenario s.

#### **Scenario-based Lookahead**

• Then we compute a sample of average approximation (SAA):

$$\tilde{J}_{k+1}(x_{k+1}) = \sum_{s=1}^{S} r_s C_s(x_{k+1})$$

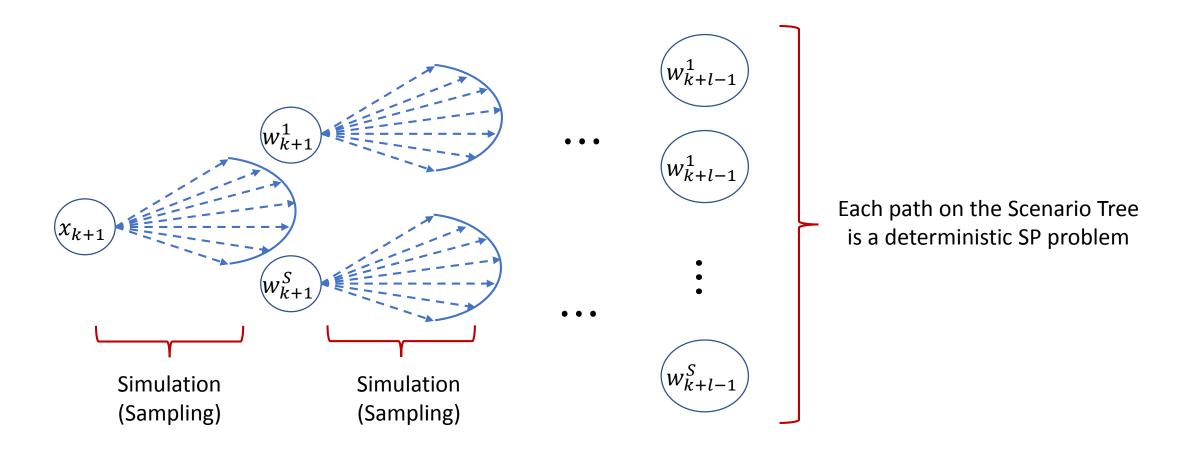
• Where  $r=(r_1,\ldots,r_S)$  are some probability distribution vector, representing the "frequency" of each scenario, often called the scenario weights.

Note that each scenario is independent.

• This idea of replacing required expectations by SAA and using simulation will be pervasive in most (if not all) algorithms in Approximate DP and Reinforcement Learning.

#### **Scenario-based Lookahead**

• The scenarios can be represented by a Scenario Tree (also called Monte-Carlo Tree):



• Let's focus on the **Deterministic** 1-step lookahead algorithm (for simplicity):

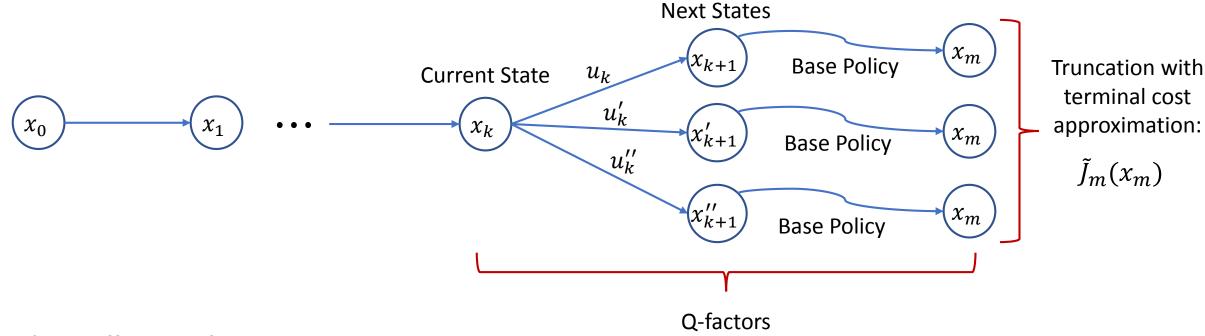
$$\tilde{\mu}_k(x_k) \in \arg\min_{u_k \in U_k(x_k)} \left\{ g_k(x_k, u_k) + \tilde{J}_{k+1}(f_k(x_k, u_k)) \right\}, \forall k \in \{0, ..., N-1\}$$

• Where the approximate function  $\tilde{J}_{k+1}(x_{k+1})$  is given as the total cost of some Base Policy (or Base Heuristic)  $\hat{\pi} = (\hat{\mu}_{k+1}, ..., \hat{\mu}_{N-1})$ :

$$\tilde{J}_{k+1}(x_{k+1}) = g_N(x_N) + \sum_{i=k+1}^{N-1} g_i(x_i, \hat{\mu}_i(x_i))$$

• Note that we can truncate the rollout at some stage m < N, and use some approximation for the terminal cost (like we did for the lookahead problems).

Schematically, we can represent the Deterministic Rollout Algorithm as follows:



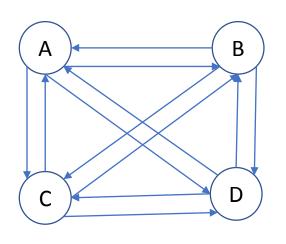
• The Rollout Policy:

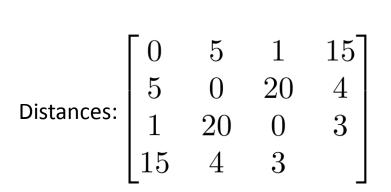
$$\tilde{\mu}_k(x_k) = \arg\min_{u_k \in U_k(x_k)} \{ \tilde{Q}_k(x_k, u_k) \}, \forall k \in \{0, ..., N-1\}$$

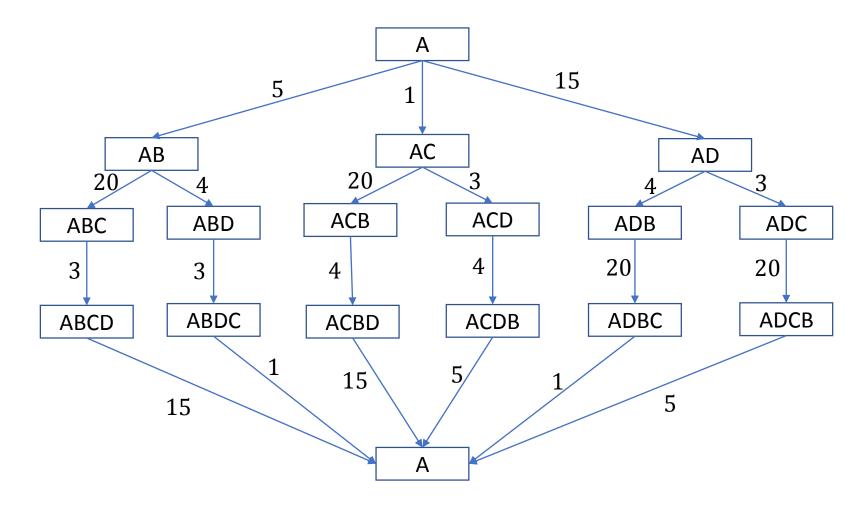
$$\tilde{Q}_k(x_k, u_k) = g_k(x_k, u_k) + \tilde{J}_{k+1}(f_k(x_k, u_k)), \forall k \in \{0, ..., N-1\}$$

# **Example: Traveling Salesman Problem**

Recall from last time the TSP:







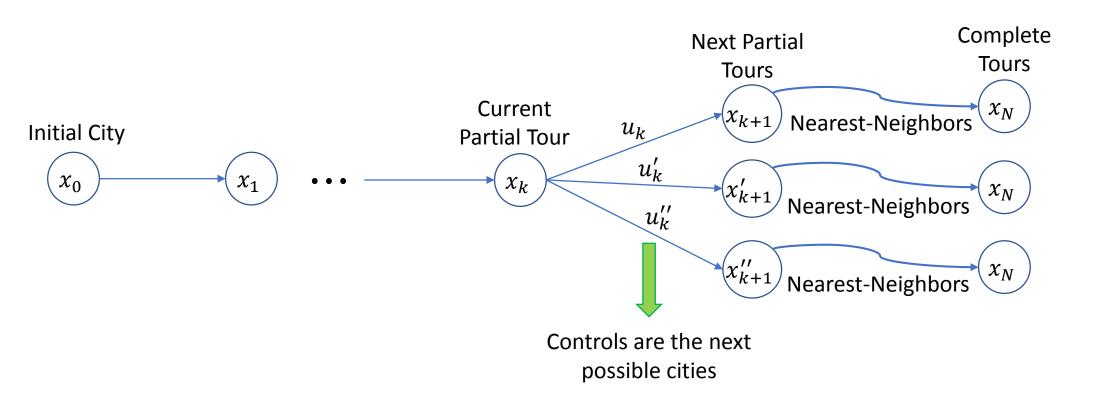
# **Example: Traveling Salesman Problem**

- The simple *Nearest-Neighbor Heuristic* can be seen as a special case of the Rollout Algorithm, where the Base Policy adds a new city to a partial tour that greedily minimize the aggregated cost and does not form a cycle.
- In our formulation  $x_k$  is a partial tour with k cities:  $x_k = (c_0, ..., c_k)$ .

• The nearest-neighbor heuristic adds a new city  $c_{k+1}$  to the partial tour if  $g(c_k, c_{k+1})$  is minimized for all cities  $c_k \neq c_0, \dots, c_k$ , Thus forming a larger partial tour:  $x_{k+1} = (c_0, \dots, c_k, c_{k+1})$ .

The Rollout Algorithm implements the nearest-neighbor heuristic as the Base Policy.

# **Example: Traveling Salesman Problem**



 After each iteration we re-run the Nearest-Neighbors heuristic from the next partial tours

- A key question in the Rollout Algorithm is whether the new rollout policy  $\tilde{\pi}$  is better than the base policy  $\hat{\pi}$ .
- This can be achieved if the Base Policy (Heuristic) is Sequentially Improving, that is:

$$\tilde{Q}_k(x_k, u_k) = g_k(x_k, u_k) + \tilde{J}_{k+1}(f_k(x_k, u_k)), \forall k \in \{0, ..., N-1\}$$
$$\tilde{J}_{k+1}(x_{k+1}) = g_N(x_N) + \sum_{i=k+1}^{N-1} g_i(x_i, \hat{\mu}_i(x_i))$$

and

$$\min_{u_k \in U_k(x_k)} \{ \tilde{Q}_k(x_k, u_k) \} \le \tilde{J}_k(x_k)$$

• In words: "The minimal approximate Q-factor at  $x_k$  is smaller then the cost-to-go of the Base Policy at  $x_k$ ".

Suppose that the Base Policy is sequentially improving. We want to show that:

$$J_{k,\tilde{\pi}}(x_k) \le \tilde{J}_k(x_k)$$

- In words: "The cost-to-go of the rollout policy at  $x_k$  is smaller than the cost-to-go of the Base Policy".
- If that inequality holds we say that the rollout policy achieves Policy Improvement.

• It turns out that if base Policy is sequentially improving, than the rollout policy achieves policy improvement.

Example: In the TSP, the nearest neighbors heuristic is sequentially improving.

• Note that the Rollout Algorithm is amenable to several base policies. Suppose we have m different base policies and at a given state  $x_{k+1}$ , the m'th base policy produces the trajectory:

$$\hat{T}_{k+1}^m = \{x_{k+1}, \hat{u}_{k+1}^m, ..., \hat{u}_{N-1}^m, \hat{x}_N^m\}$$

- And corresponding cost:  $C(\hat{T}_{k+1}^m)$
- And the approximate function  $\tilde{J}_{k+1}(x_{k+1})$  is given by:

$$\tilde{J}_{k+1}(x_{k+1}) = \min_{m} \{C(\hat{T}_{k+1}^m)\}$$

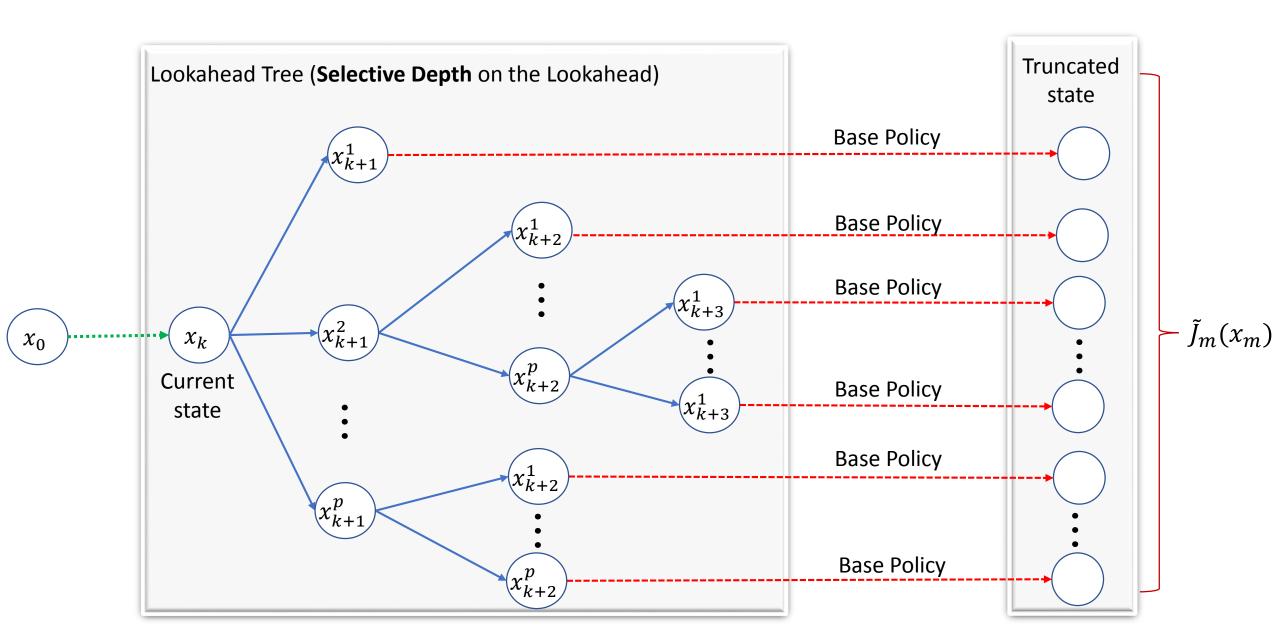
• And it follows that if every base policy is sequentially improving than the *superheuristic*, that produces the cost  $\tilde{J}_{k+1}(x_{k+1})$  given  $x_{k+1}$  is also sequentially improving.

#### **Deterministic Tree Search**

- Now let's combine everything together!
  - Multistep lookahead
  - Rollout Policy
  - Truncation with terminal cost approximation

- The idea is to use a multistep lookahead model where we selectively use base policies to provide approximations for the value functions.
- If the base policies are sequentially improving, then the policy that comes out of the lookahead problem equipped with the rollout policy will achieve policy improvement.
- This procedure leads to a *Tree Search Algorithm*, which is a powerful approximation method (specially when the base policies are Deep Neural Networks!)

#### **Deterministic Tree Search**



#### **Deterministic Tree Search**

- This formulation already exposes a central idea in approximate DP:
  - Exploration vs Exploitation
  - The idea of the Tree Search algorithm is to perform a selective depth lookahead
  - Pruning vs Expansion of the tree nodes.
  - For example, we can prune a node at some state  $x_{k+1}$  is some score function  $S_{k+1}(x_{k+1})$  falls below some threshold  $\alpha_{k+1}$ .
  - We can use the 1-step lookahead problem equipped with the Base Policy to provide the scoring function:

$$S_{k+1}(x_{k+1}) = \min_{u_{k+1} \in U_{k+1}(x_k+1)} \left\{ g_{k+1}(x_{k+1}, u_{k+1}) + \tilde{J}_{k+2}(f_{k+1}(x_{k+1}, u_{k+1})) \right\}$$
$$\tilde{J}_{k+2}(x_{k+2}) = \tilde{J}_m(x_m) + \sum_{i=k+2}^{m-1} g_i(x_i, \hat{\mu}_i(x_i))$$

## **Rollout Algorithm: Stochastic Case**

 On the stochastic case the notion of policy improvement carries out nicely, and we still have:

$$J_{k,\tilde{\pi}}(x_k) \le \tilde{J}_k(x_k)$$

• But the key question now is how to execute the Base Policy  $\hat{\pi}$  in order to obtain the rollout trajectories.

• We do so by using simulation/sampling. Namely starting from  $x_k$ , we compute:

$$x_{i+1} = f_i(x_i, \hat{\mu}_i(x_i), w_i), \forall i = k+1, ..., N-1$$

• For sampled sequences  $(w_k, ..., w_{N-1})$ .

## **Rollout Algorithm: Stochastic Case**

• We can use SAA, to compute the expected Q-factors as follows:

$$\tilde{Q}_{k}(x_{k}, u_{k}) = \mathbb{E}_{w_{k}}[g_{k}(x_{k}, u_{k}, w_{k}) + \tilde{J}_{k+1}(f_{k}(x_{k}, \hat{\mu}_{k}(x_{k}), w_{k}))]$$

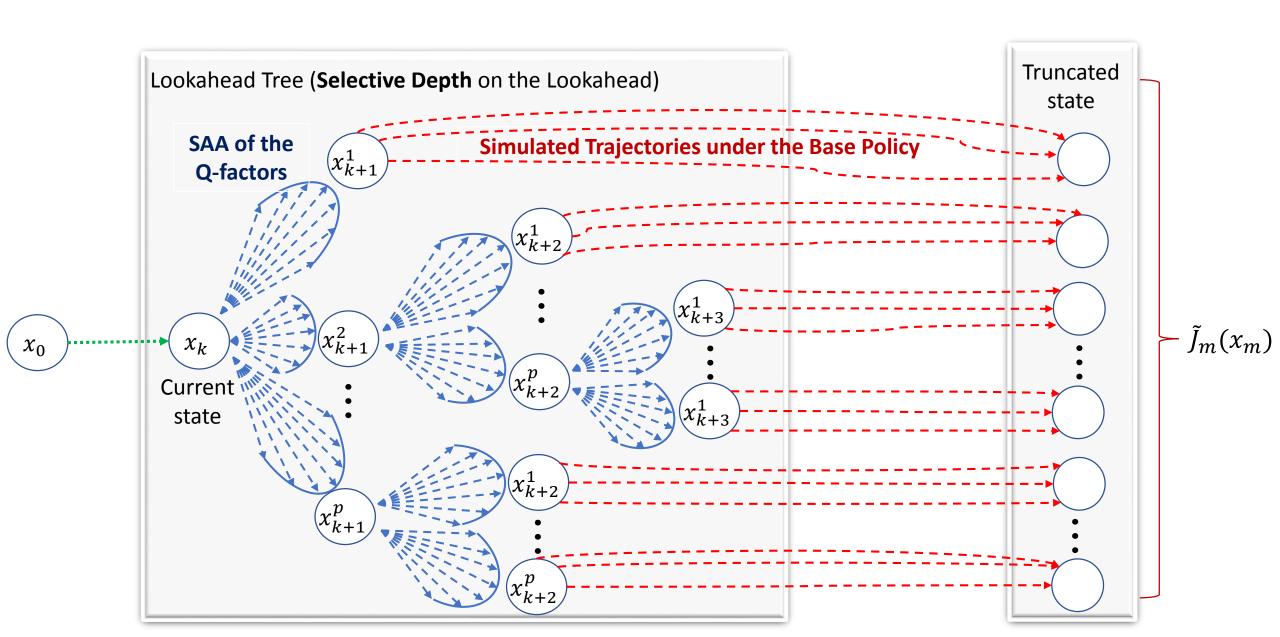
$$\tilde{Q}_{k}(x_{k}, u_{k}) \approx \sum_{s=1}^{S} r_{s}(g_{k}(x_{k}, u_{k}, w_{k}^{s}) + \tilde{J}_{k+1}(f_{k}(x_{k}, \hat{\mu}_{k}(x_{k}), w_{k}^{s})))$$

Where S is the number of samples.

• Then the **rollout policy** becomes:

$$\tilde{\mu}_k(x_k) \in \min_{u_k \in U_k(x_k)} \left\{ \sum_{s=1}^S r_s \left( g_k(x_k, u_k, w_k^s) + \tilde{J}_{k+1}(f_k(x_k, \hat{\mu}_k(x_k), w_k^s)) \right) \right\}$$

## Monte-Carlo Tree Search (MCTS)



## **Monte-Carlo Tree Search (MCTS)**

- The Monte-Carlo Tree Search is a very powerful and flexible framework to perform Approximations in the Value Space.
- With the implementation of such schemes, several questions can be raised:
  - (1) How can we score good control options now? **Exploration vs Exploitation**
  - (2) How to obtain reasonable terminal costs at the truncation states?
  - (3) How can we efficiently obtain samples?
  - (4) How to avoid performing so much simulation online?
  - (5) How can we "control" the simulation error? What about the Variance?
  - (6) What would be a good Base Policy?

• Successful cases, such as AlphaGo, provide answers to the above questions. We will continue next lecture, with potential answers!