## An introduction to the theory of SDDP algorithm

V. Leclère (ENPC)

October 29, 2013

### Introduction

- Large scale stochastic problem are hard to solve.
- Two ways of attacking such problems :
  - decompose (spatially) the problem and coordinate solutions,
  - construct easily solvable approximations (Linear Programming).
- Behind the name SDDP there is three different things:
  - a class of algorithm,
  - a specific implementation of the algorithm,
  - a software implementing this method develloped by PSR
- The aim of this talk is to give you an idea of how the class of algorithm is working.

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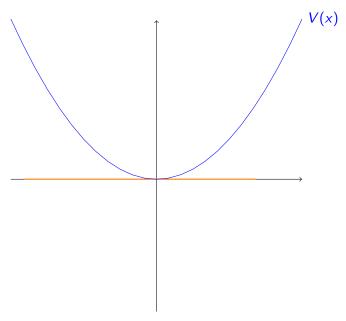
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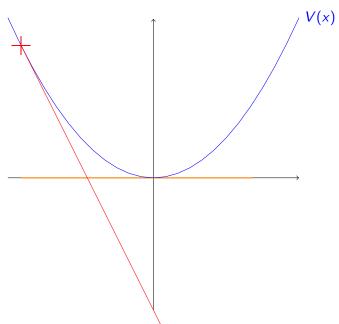
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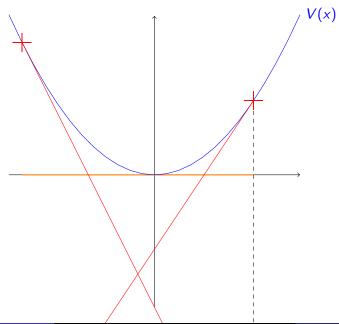
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  - Problem statement
  - Some background on Dynamic Programming
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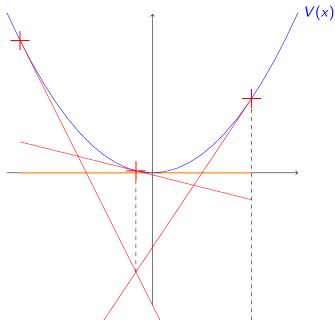
Kelley's algorithm

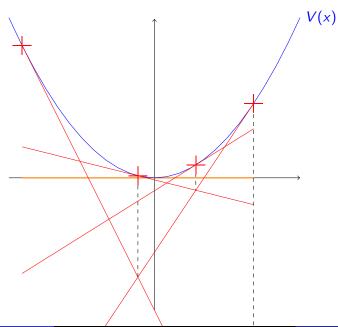


Kelley's algorithm









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### Problem considered

We consider a discrete and finite time optimal control problem

$$\min_{u \in \mathbb{U}^{T}} \sum_{t=0}^{T-1} L_{t}(x_{t}, u_{t}) + K(x_{T}),$$

$$s.t. \quad x_{t+1} = f_{t}(x_{t}, u_{t}).$$

- Where
  - $x_t \in \mathbb{X}$  is the state at time t,
  - $u_t \in \mathbb{U}$  the control applied at time t.
- We assume that
  - f<sub>t</sub> are linear.
  - ullet U and X are compact.
- We consider convex cost  $L_t(x_t, u_t)$ , and a final cost  $K(x_T)$ .
- A policy is a sequence of functions  $\pi = (\pi_1, \dots, \pi_{T-1})$  giving for any state x a control u.

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### Introducing Bellman's function

This problem can be solved by dynamic programming. In this case we introduce the Bellman function defined by

$$\begin{cases} V_T(x) &= K(x), \\ V_t(x) &= \min_{u_t \in \mathbb{U}} \left\{ L_t(x, u_t) + V_{t+1} \circ f_t(x, u_t) \right\} = \mathcal{T}_t(V_{t+1})(x) \end{cases}$$

where

$$\mathcal{T}_t(A): x \mapsto \min_{u_t \in \mathbb{U}} \left\{ L_t(x, u_t) + A \circ f_t(x, u_t) \right\}.$$

$$\pi_t(x) \in \operatorname*{arg\,min}_{u_t \in \mathbb{U}} \left\{ L_t(x, u_t) + V_{t+1} \circ f_t(x, u_t) \right\}$$

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Indeed an optimal policy for this problem is given by

$$\pi_t(x) \in \arg\min_{u_t \in \mathbb{T}} \left\{ L_t(x, u_t) + V_{t+1} \circ f_t(x, u_t) \right\}$$

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### Properties of Bellman operator

• Monotonicity:

$$\forall x \in \mathbb{X}, \quad V(x) \leq \overline{V}(x) \quad \Rightarrow \quad \forall x \in \mathbb{X}, \quad (\mathcal{T}V)(x) \leq (\mathcal{T}\overline{V})(x).$$

• Convexity: if  $L_t$  is jointly convex in (x, u), V is convex, and  $f_t$  is affine then

$$x \mapsto (\mathcal{T}V)(x)$$
 is convex.

• Linearity: for any piecewise linear function V, if  $L_t$  is also piecewise linear, and  $f_t$  affine, then

$$x \mapsto (\mathcal{T}V)(x)$$
 is piecewise linear.

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### **Duality property**

- Consider  $J: \mathbb{X} \times \mathbb{U} \to \mathbb{R}$  jointly convex.
- Define

$$\varphi(x) = \min_{u \in \mathbb{U}} J(x, u),$$

• Then we can obtain a subgradient  $\lambda \in \partial \varphi(x_0)$  as the dual multiplier of

$$\min_{x,u} J(x,u),$$
s.t.  $x_0 - x = 0$   $[\lambda]$ 

(This is the marginal interpretation of the multiplier).

In particular it means that

$$\varphi(\cdot) > \varphi(x_0) + \langle \lambda, \cdot - x_0 \rangle.$$

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### General idea

- The SDDP algorithm recursively constructs an approximation of each Bellman function as the supremum of a number of affine functions.
- At stage k we have  $V_t^{(k)}$  lower approximations of  $V_t$  and we want to construct a better approximation.
- We follow an optimal trajectory  $(x_t^{(k)})_t$  of the approximated problem and add a cut for each Bellman function.

# Stage k of SDDP description (1/2)

- Began a loop forward in time by setting t = 0 and  $x_t^{(k)} = x_0$ ,
- solve

$$\min_{x,u} L_t(x,u) + V_{t+1}^{(k)} \circ f_t(x,u),$$

$$x = x_t^{(k)}. [\lambda_t^{(k+1)}]$$

- We call
  - $\beta_t^{(k+1)}$  the value of the problem,
  - $\lambda_t^{(k+1)}$  a multiplier of the constraint  $x = x_t^{(k)}$ ,
  - $u_t^{(k)}$  an optimal control.
- This can also be written as

$$\beta_t^{(k+1)} = \mathcal{T}_t \left( V_{t+1}^{(k)} \right) \left( x_t^{(k)} \right),$$
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# Stage k of SDDP description (2/2)

Thus,

$$\beta_{t}^{(k+1)} + \langle \lambda_{t}^{(k+1)}, \cdot - x_{t}^{(k)} \rangle \leq \mathcal{T}_{t} \left( V_{t+1}^{(k)} \right) \leq \mathcal{T}_{t} \left( V_{t+1} \right) = V_{t}.$$

- Thus  $x \mapsto \beta_t^{(k+1)} + \left\langle \lambda_t^{(k+1)}, x x_t^{(k)} \right\rangle$  is a cut.
- ullet We update our approximation of  $V_t$  by defining

$$V_t^{(k+1)} = \max\left\{V_t^{(k)}, \beta_t^{(k+1)} + \left\langle\lambda_t^{(k+1)}, \cdot - x_t^{(k)}\right\rangle\right\}.$$

- $V_t^{(k+1)}$  is convex and lower than  $V_t$ .
- set

$$x_{t+1}^{(k)} = f_t \left( x_t^{(k)}, u_t^{(k)} \right)$$

• Upon reaching time t = T we have completed iteration k of the algorithm.

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Conclusion

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### Initialisation and stopping rule

- To initialize the algorithm it seems that we need a lower bound (that exist) to all value function.
- In fact we can choose  $V_t^{(0)} = 0$  in order to compute the cuts, and simply set  $V_t^{(1)}$  equal to the first cut, which means that we "forget"  $V_t^{(0)}$  in the maximum that determine  $V_t^{(1)}$ .
- At any step k we have a admissible, non optimal solution  $(u^{(k)})_t$ , with
  - an upper bound

$$\sum_{t=0}^{T-1} L_t \left( x_t^{(k)}, u_t^{(k)} \right) + K \left( x_T^{(k)} \right),$$

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- a lower bound  $V_0^{(k)}(x_0)$ .
- A reasonable stopping rule for the algorithm is given by checking that the (relative) difference of the upper and lower bound is small.

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### What's new?

Now we introduce some random variables  $\mathbf{W}_t$  in our problem. This complexify the algorithm in different ways :

- we need some probabilistic assumptions;
- for each stage k we need to do a forward phase that yields a trajectory  $(x_t^{(k)})_t$ , and a backward phase that gives a new cut;
- we can not compute an exact upper bound for the problem's value.

### Problem statement

$$\begin{aligned} & \underset{\pi}{\text{min}} & & \mathbb{E}\left(\sum_{t=0}^{T-1} L_t(\mathbf{X}_t, \mathbf{U}_t, \mathbf{W}_t) + K(\mathbf{X}_T)\right), \\ & s.t. & & \mathbf{X}_{t+1} = f_t(\mathbf{X}_t, \mathbf{U}_t, \mathbf{W}_t), \\ & & & \mathbf{U}_t = \pi_t(\mathbf{X}_t, \mathbf{W}_t). \end{aligned}$$

Stochastic case

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Where  $(\mathbf{W}_t)_{t \in \{1,\dots,T\}}$  is assumed to be a white noise.

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## Stochastic Dynamic Programming

This problem can be solved by dynamic programming. In this case we introduce the Bellman function defined by

$$\begin{cases}
V_{T}(x) &= K(x), \\
\hat{V}_{t}(x, w) &= \min_{u_{t} \in \mathbb{U}} L_{t}(x, u_{t}, w) + V_{t+1} \circ f_{t}(x, u_{t}, w), \\
V_{t}(x) &= \mathbb{E} \left( \hat{V}_{t}(x, \mathbf{W}_{t}) \right).
\end{cases} (1)$$

Indeed an optimal policy for this problem is given by

$$\pi_t(x, w) \in \operatorname*{arg\,min}_{u_t \in \mathbb{I}} \left\{ L_t(x, u_t, w) + V_{t+1} \circ f_t(x, u_t, w) \right\}$$

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### Bellman operator

For any time t, and any function A mapping the set of states and noises  $\mathbb{X} \times \mathbb{W}$  into  $\mathbb{R}$  we define :

$$\hat{\mathcal{T}}_t(A)(x,w) := \min_{u_t \in \mathbb{U}} L_t(x,u_t,w) + A \circ f_t(x,u_t,w).$$

Thus the Bellman equation simply reads

$$\begin{cases} V_{\mathcal{T}}(x) &= K(x), \\ V_{t}(x) &= \mathcal{T}_{t}(V_{t+1})(x) := \mathbb{E}\left(\hat{\mathcal{T}}_{t}(V_{t+1})(x, \mathbf{W}_{t})\right). \end{cases}$$

The Bellman operator have the same properties as in the deterministic case.

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## Duality theory (1/2)

Consider that we know  $V_{t+1}^{k+1} \leq V_{t+1}$ .

$$\hat{\beta}_{t}^{(k+1)}(w) = \min_{x,u} \quad L_{t}(x,u,w) + V_{t+1}^{(k+1)} \circ f_{t}(x,u,w),$$

$$s.t \quad x = x_{t}^{(k)} \qquad [\hat{\lambda}_{t}^{(k+1)}(w)]$$

Which can also be written

$$\hat{\beta}_t^{(k+1)}(w) = \hat{\mathcal{T}}_t \left( V_{t+1}^{(k)} \right) (x, w),$$
$$\hat{\lambda}_t^{(k+1)}(w) \in \partial_x \hat{\mathcal{T}}_t \left( V_{t+1}^{(k)} \right) (x, w)$$

Thus for all w.

$$\hat{\beta}_t^{(k+1)}(w) + \left\langle \hat{\lambda}_t^{(k+1)}(w), x - x_t^{(k)} \right\rangle \leq \hat{\mathcal{T}}_t \left( V_{t+1}^{(k)} \right) (x, w) \leq \hat{V}_t(x, w).$$

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# Duality theory (2/2)

Thus we have an affine minorant for each realisation of  $\mathbf{W}_t$ . Replacing w by the random variable  $\mathbf{W}_t$  and taking the expectation yields the following affine minorant

$$\beta_t^{(k+1)} + \left\langle \lambda_t^{(k+1)}, \cdot - x_t^{(k)} \right\rangle \leq V_t,$$

where

$$\left\{ \begin{array}{ll} \boldsymbol{\beta}_t^{(k+1)} & := & \mathbb{E}\left(\hat{\boldsymbol{\beta}}_t^{(k+1)}(\mathbf{W}_t)\right) = \mathcal{T}_t\left(\boldsymbol{V}_{t+1}^{(k)}\right)(\boldsymbol{x}), \\ \boldsymbol{\lambda}_t^{(k+1)} & := & \mathbb{E}\left(\hat{\boldsymbol{\lambda}}_t^{(k+1)}(\mathbf{W}_t)\right) \in \partial_{\boldsymbol{x}}\mathcal{T}_t\left(\boldsymbol{V}_{t+1}^{(k)}\right)(\boldsymbol{x}). \end{array} \right.$$

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## At the beginning of step

At the beginning of step k we suppose that we have, for each time step t an approximation  $V_t^k$  of  $V_t$  verifying

- $V_t^k \leq V_t$ ,
- $V_T^k = K,$
- $V_t^k$  is convex.

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## Forward path: define a trajectory

- Randomly select a scenario  $(w_0, \ldots, w_{T-1}) \in \mathbb{W}^T$ .
- Define a trajectory  $(x_t^{(k)})_{t=0,\dots,T}$  by

$$x_{t+1}^{(k)} = f_t(x_t^{(k)}, u_t^{(k)}, w_t),$$

where  $u_t^{(k)}$  is an optimal solution of

$$\min_{u\in\mathbb{U}}L_t\left(x_t^{(k)},u,w_t\right)+V_{t+1}^{(k)}\circ f_t\left(x_t^{(k)},u,w_t\right).$$

• This trajectory is given by the optimal policy where  $V_t$  is replaced by  $V_t^{(k)}$ .

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### Backward path: add cuts

- ullet For any t we want to add a cut to the approximation of  $V_t$ .
- At time t solve, for any w possible

$$\hat{\beta}_{t}^{(k+1)}(w) = \min_{x,u} \quad L_{t}(x,u,w) + V_{t+1}^{(k+1)} \circ f_{t}(x,u,w),$$

$$s.t \quad x = x_{t}^{(k)} \qquad [\hat{\lambda}_{t}^{(k+1)}(w)]$$

- $\begin{array}{l} \bullet \ \ \mathsf{Compute} \ \lambda_t^{(k+1)} = \mathbb{E} \left( \lambda_t^{(k+1)}(\mathbf{W}_t) \right) \ \mathsf{and} \\ \beta_t^{(k+1)} = \mathbb{E} \left( \beta_t^{(k+1)}(\mathbf{W}_t) \right). \end{array}$
- Add a cut

$$V_t^{(k+1)}(x) = \max\left\{V_t^{(k)}(x), \beta_t^{(k+1)} + \left\langle \lambda_t^{(k+1)}, x - x_t^{(k)} \right\rangle\right\}$$

• Go one step back in time :  $t \leftarrow t - 1$ . Upon reaching t = 0 we have completed step k of the algorithm.

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### Initialization and stopping rule

- In order to accelerate the convergence it can be useful to bypass a few forward paths by abritrarily choosing some trajectories  $(x_t^{(k)})_t$ .
- We have a lower bound given by  $V_0^{(k)}(x_0)$ .
- The upper bound is more complicated (expectation over the whole process  $(W_0, \ldots, W_{T-1})$ , but can be estimated by Monte-Carlo methods, and we have no control over the error of our solution.
- A heuristic stopping rule consist in stopping the algorithm if the lower bound is in the confidence interval of the upper bound for a determined number of Monte-Carlo simulation.

### A few other implementation

- We presented DOASA : select one scenario (one realisation of  $(W_1, \ldots, W_{T-1})$ ) to do a forward and backward path.
- Classical SDDP: select a number N of scenarios to do the forward path (computation can be parallelized). Then during the backward path we add N cuts to  $V_t$  before computing the cuts on  $V_{t-1}$ .
- CUPPS algorithm suggest to use  $V_{t+1}^{(k)}$  instead of  $V_{t+1}^{(k+1)}$  in the computation of the cuts. In practice :
  - select randomly a scenario  $(w_t)_{t=0,...,T-1}$ ;
  - at time t we have a state  $x_t^{(k)}$ , we compute the new cut for  $V_t$ ;
  - choose the optimal control corresponding to the realization  $W_t = w_t$  in order to compute the state  $x_{t+1}^{(k)}$  where the cut for  $V_{t+1}$  will be computed, and goes to the next step.
- We can compute some cuts before starting the algorithm. For example by bypassing the forward phase by choosing the trajectory  $(x_t^{(k)})_{t=0,\dots,T}$ .

Stochastic case

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- The problem studied was risk neutral.
- However a lot of works has been done recently about how to solve risk averse problems.
- Most of them are using CVAR, or a mix between CVAR and expectation.
- Indeed CVAR can be used in a linear framework by adding another variable.
- Another easy way is to use "composed risk measures".
- Finally a convergence proof with convex costs (instead of linear costs) exists. However it require to solve non-linear problems.

Conclusion

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#### Conclusion

SDDP is an algorithm, more precisely a class of algorithms that

- exploit convexity of the value functions (from convexity of costs...);
- does not require discretization;
- construct outer approximations of  $V_t$ , those approximations being precise only "in the right places";
- gives bounds :
  - real lower bound  $V_0^{(k)}(x_0)$ ,
  - estimated (by Monte-Carlo) upper bound;
- construct linear-convex approximations, thus enabling to use linear solver like CPLEX,
- have some proof of asymptotic convergence.





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