

Optimizing the exercise boundary for the holder of an American option over a parametric family

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American options

The holder has the right to exercise the option at any time $t \in [0, T]$ and get a payoff $V(X_t, t)$, where the coordinates of $X_t \in \mathbb{R}^d$ are the prices of some underlying assets at time t . The **diffusive process** X_t solves

$$dX = a(X)dt + \sigma(X)dW \quad (1)$$

with **generator** $\mathcal{L} = \nabla \cdot a + \frac{1}{2}\text{tr}(\sigma\sigma^T D^2)$ where D^2 is the Hessian. The optimization problem is to find

$$\sup_{\tau \in \mathcal{S}_{0,T}} E[V(X_\tau, \tau)], \quad (2)$$

where $\mathcal{S}_{0,T}$ is the set of all **stopping times** bounded by the final time T .

Let $w(x, t)$ be the solution of the optimization problem starting with prices x at the time t . Clearly, $w \geq V$, and if we denote by Ω the points (x, t) for which $w > V$, then w solves the parabolic PDE

$$w_t + \mathcal{L}w = 0, \quad \text{when } (x, t) \in \Omega \quad (3)$$

with boundary conditions

$$\begin{aligned} w(x, t) &= V(x, t), & \text{for } (x, t) \in \partial\Omega \text{ or } t = T, \\ \nabla w(x, t) &= \nabla V(x, t), & \text{for } (x, t) \in \partial\Omega. \end{aligned} \quad (4)$$

The boundary conditions (4) determine $\partial\Omega$.

① Finite differences

- Works up to $d \simeq 4$.
- Computational work has exponential growth with d .
- Computational memory has exponential growth with d .

② Longstaff and Schwartz.

- Works up to $d \simeq 10$.
- It is a Monte Carlo method, but the confidence intervals have exponential growth with d .

Our objective is to design a Monte Carlo method such that the confidence intervals have polynomial growth with d .

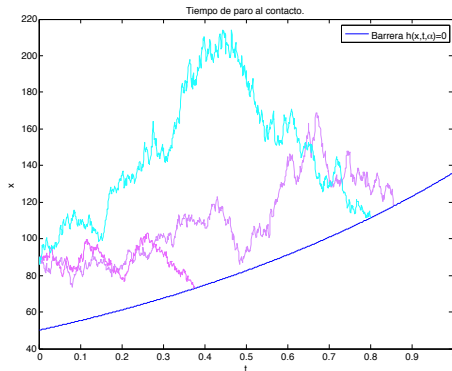
Parametrized free boundary

We optimize only over stopping times given by a level curve.

$$\tau_\theta = \inf \{t \geq 0 \mid h(X_t, t, \theta) \leq 0\} . \quad (5)$$

Our objective is to maximize the function

$$f(\theta) = E[V(X_{\tau_\theta}, \tau_\theta)] .$$



We have random variables V_θ , indexed by the parameter θ . The aim is to maximize

$$f(\theta) = E[V_\theta]. \quad (6)$$

Difficulty:

- $f(\theta)$ is evaluated by the Monte Carlo method, which means that evaluations are expensive and have a significant statistical error.

$$\hat{f}(\theta) = \frac{1}{L} \sum_{k=1}^L V_k \approx f(\theta) + \frac{\sigma_\theta}{\sqrt{L}} Z, \quad Z \sim N(0, 1). \quad (7)$$

Solution: Find a random variable W_θ that satisfies

$$\nabla f(\theta) \approx E[W_\theta]. \quad (8)$$

Desiderata:

- Low *bias*,

$$\text{bias} = E[W_\theta] - \nabla f(\theta).$$

- Evaluation of W_θ should be as hard as evaluating V_θ .
- The standard deviations of W_θ and V_θ should be of the same order.

In general, there is **no rule** for the construction of W_θ . The construction in each application is the **sensitivity analysis**.

Robbins-Monro optimization

A general iteration takes the form

$$\hat{\theta}_{n+1} = \hat{\theta}_n - \frac{1}{n^\gamma} K_n W_n, \quad (9)$$

- *Stochastic Gradient Descent*: $K_n = \lambda I$, $\gamma = 1.0$
- *Optimal*: $K_n \approx H^{-1}$ where H is the Hessian of f evaluated at α_* .
- Big difference when H is an **ill conditioned** matrix
- Since W_{θ_n} has mean $\nabla f(\theta_n)$, the slope of a least square fit of the points (θ_n, W_{θ_n}) will have a slope approximately equal to the derivative of the map $\theta \rightarrow \nabla f(\theta)$; this is, the Hessian of f .

Linear regression approximation of H^{-1}

Let X_n and Y_n be the corresponding $n \times (p + 1)$ and $n \times p$ matrices with entries

$$x_k = (\hat{\theta}_k, 1), \quad y_k = W_k,$$

consider the linear regression $Y = XB$ and denote the first p rows of a matrix M by \overline{M} . We calculate the natural estimators

$$\begin{aligned} B_n &= (X_n^T X_n)^{-1} X_n^T Y, & H_n &= \overline{B_n} \\ G_n &= \overline{B_n}^{-1}, & K_n &= \frac{G_n + G_n^T}{2}, \end{aligned} \tag{10}$$

and use K_n as our H^{-1} estimator and $\gamma = 0.6$.

- The estimator is $\overline{\theta}_n = \frac{1}{n}(\hat{\theta}_1 + \dots + \hat{\theta}_n)$, **Polyak averaging**.
- Similar methods (with $\gamma = 1$ and no Polyak averaging) were analyzed by **Lai and Robbins in 1981**, without numerical simulations.
- The resulting optimization is **Affine Invariant**:

$$B \theta_O(f \circ B) = \theta_O(f), \quad B : \mathcal{S} \longrightarrow \mathcal{S} \quad \text{linear}. \tag{11}$$

PDE representation (Feynman-Kac).

- The function $u(x, t, \theta) = E_{x,t}[V(X_{\tau_\theta}, \tau_\theta)]$ satisfies the PDE

$$u_t + \mathcal{L}u = 0, \quad \text{cuando } h(x, t, \theta) > 0 \quad (12)$$

with boundary conditions

$$u(x, t, \theta) = V(x, t), \quad \text{cuando } h(x, t, \theta) = 0 \text{ o } t = T. \quad (13)$$

- The function $\dot{u}(x, t, \theta) = \nabla_\theta u(x, t, \theta)$ satisfies the same PDE but with boundary conditions

$$\dot{u}(x, t, \theta) = \left(\frac{\partial u}{\partial n} - \frac{\partial V}{\partial n} \right) \frac{\dot{h}}{\|\nabla h\|} \quad \text{cuando } h(x, t, \theta) = 0 \quad (14)$$

$$\dot{u}(x, T, \theta) = 0, \quad (15)$$

The vector $n = \frac{\nabla h}{\|\nabla h\|}$ is the **inner normal** of the hypersurface $h(x, t, \theta) = 0$ in R^d .

Probabilistic representation of \dot{u}

The Feynman-Kac formula gives

$$\nabla f(\theta) = \dot{u}(X_0, 0, \theta) = E_{X_0, 0} \left[\left(\frac{\partial u}{\partial n} - \frac{\partial V}{\partial n} \right) \frac{\dot{h}}{\|\nabla h\|} \right]. \quad (16)$$

All functions are evaluated at $(X_{\tau_\theta}, \tau_\theta, \theta)$. It is tempting to define

$$W_\theta = \left(\frac{\partial u}{\partial n} - \frac{\partial V}{\partial n} \right) \frac{\dot{h}}{\|\nabla h\|}$$

There is an obvious obstacle: **We need to evaluate $\frac{\partial u}{\partial n}$ at the boundary.**

We introduce a simulation algorithm for diffusion processes (time increment Δt) which is very efficient close to the boundary, and therefore, useful when estimating $\frac{\partial u}{\partial n}$. The (X_τ, τ) simulation will be denoted by $(\hat{X}_{\hat{\tau}}, \hat{\tau})$. **Main properties:**

- $E[V(\hat{X}_{\hat{\tau}}, \hat{\tau})] = E_{x,t}[V(X_\tau, \tau)] + O(\varepsilon \Delta t^\gamma)$
- $\text{Var } V(\hat{X}_{\hat{\tau}}, \hat{\tau}) = O(\varepsilon)$
- $E[\hat{\tau}] = O(\varepsilon)$

where ε is the distance to the boundary $h = 0$.

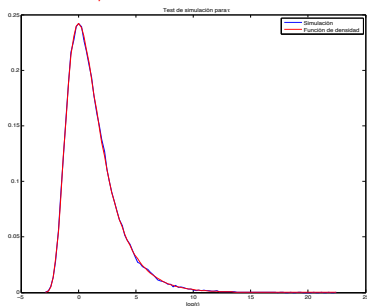
We have $\gamma = 1$ when $d = 1$ and $\gamma = 1/2$ in high dimensions ($d > 1$). **The bias, variance and simulation time are $O(\varepsilon)$.**

Stopping time simulation.

Assume $\bar{X}_t \in \mathbb{R}$ is a Brownian motion and the linearized boundary is $\bar{h}(x, t) = c - x$, this is, the line $x = c > 0$. The density function for $\bar{\tau}$ is given by

$$P(\bar{\tau} \in t + dt) = \frac{1}{\sqrt{2\pi t^3}} c e^{-c^2/2t} dt + o(dt). \quad (17)$$

Changing variables with $z = c/\sqrt{t}$ we obtain $\bar{\tau} = c^2/Z^2$ with $Z \sim N(0, 1)$.



Since **paths started close to the boundary do not last long**, we can generate many paths and take an average:

$$\hat{u} = \varepsilon \sum_{i=1}^{1/\varepsilon} V(\hat{X}_{\hat{\tau}_i}, \hat{\tau}_i), \quad (18)$$

and we obtain

- $E[\hat{u}] = E_{x,t}[V(X_\tau, \tau)] + O(\varepsilon \Delta t^\gamma)$
- $\text{Var } \hat{u} = O(\varepsilon^2)$
- $E[\text{Work}(\hat{u})] = \frac{1}{\varepsilon} \frac{E[\hat{\tau}]}{\Delta t} = O\left(\frac{1}{\Delta t}\right)$

The time it takes to generate $\frac{1}{\varepsilon}$ paths started at distance ε from the boundary is the same time it takes to generate one path far from the boundary. We call this technique the **Splash Method**.

The estimator for the normal derivative at a point x at the boundary $h = 0$ is of course

$$D = \lim_{\epsilon \rightarrow 0} \frac{\hat{u}(x + \epsilon n) - V(x, 0)}{\epsilon} .$$

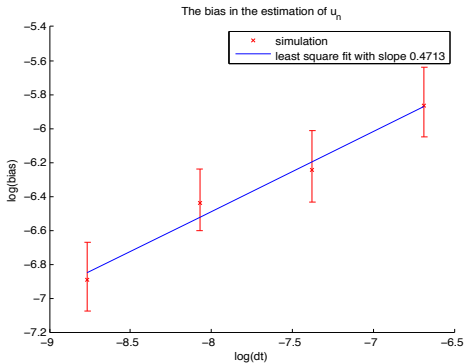
The properties of the estimator \hat{u} give

- $E[D] = \partial_n u + O(\Delta t^\gamma)$
- $\text{Var } D = O(1)$

The only task left is to evaluate the limit as $\epsilon \rightarrow 0$ in the definition of D .

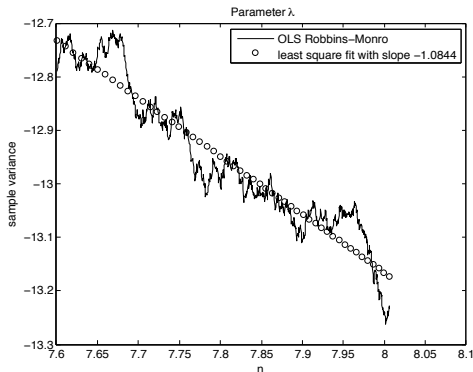
Numerical simulation of D , $d = 3$

- $d \ln(X) = \sigma dW$
- $u(x, t) = \exp\left(-\frac{1}{2}(T - t)\xi^T(\sigma \sigma^T) \xi\right) \cos(\xi^T \ln(x))$.
- $h(x, t) = \left(1 + \frac{t}{2}\right)^2 - \|x - c\|^2$.

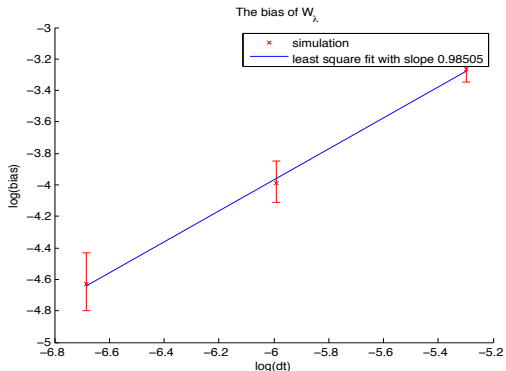


Numerical simulations. $d = 1$ and two parameters.

- $dX = r X dt + \sigma X dW$ Risk neutral process
- $h(x, t, S, \lambda) = x - S e^{\lambda t}$ Exponential boundary
- $V(x, t) = e^{-r t} (K - x)_+$ American put option
- We can calculate $u(x, t, S, \lambda)$ analytically.
- Convergence speed is measured by $\text{Var}(S_n)$ y $\text{Var}(\lambda_n)$.

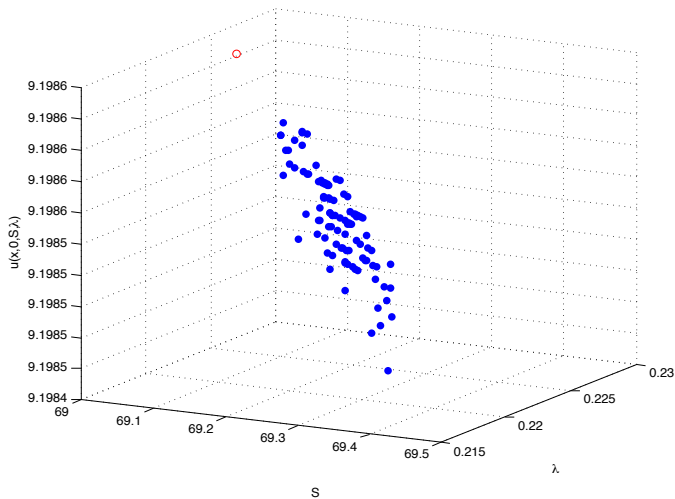


- bias = $\nabla f(\theta) - E[W_\theta]$
- The bias is $O(\Delta t)$ for $d = 1$ and $O(\sqrt{\Delta t})$ for $d > 1$.



100 simulations of 3000 steps

The price of the put option with the simulation parameters is **\$9.22** .



- José Vidal Alcalá Burgos, Optimizing the exercise boundary for the holder of an American option over a parametric family, Ph.D. Thesis, ProQuest (2012).
<http://gradworks.umi.com/35/24/3524127.html>
- Code for the *affine invariant optimization* is available at <https://github.com/vidalalcala/sopt-ols>

Efficient online update

There is an **online update** algorithm for K_n , the estimator of H^{-1} .

$$\begin{aligned} s_{n+1} &= \frac{1}{1 + x_{n+1} P_n x_{n+1}^T} \\ u_{n+1} &= s_{n+1} \overline{P_n x_{n+1}^T} \\ v_{n+1} &= y_{n+1} - x_{n+1} \overline{B_n} \\ t_{n+1} &= \frac{1}{1 + v_{n+1} G_n u_{n+1}} \\ G_{n+1} &= G_n - t_{n+1} G_n u_{n+1} v_{n+1} G_n \\ \overline{B_{n+1}} &= \overline{B_n} + s_{n+1} P_n x_{n+1}^T v_{n+1} \\ P_{n+1} &= P_n - s_{n+1} P_n x_{n+1}^T x_n P_n^T \end{aligned} \tag{19}$$

- $P_n = (X_n^T X_n)^{-1}$ is the **precision matrix**.
- $O(p^2)$ **paralellizable** operations.

Remove bias , $\varepsilon \rightarrow 0$.

$$E_\varepsilon[V(\hat{X}_{\hat{\tau}}, \hat{\tau})] - V(x, 0) = E_\varepsilon[V(\hat{X}_{\hat{\tau}}, \hat{\tau}) - V(x, 0) \mid \hat{\tau} \geq \Delta t]P_\varepsilon(\hat{\tau} \geq \Delta t) \\ + E_\varepsilon[(V(\hat{X}_{\hat{\tau}}, \hat{\tau}) - V(x, 0)) 1_{\hat{\tau} < \Delta t}]$$

- $E_\varepsilon[V(\hat{X}_{\hat{\tau}}, \hat{\tau}) - V(x, 0) \mid \hat{\tau} \geq \Delta t] = O(1)$, $S(y) = \frac{y}{\Delta t} e^{-y^2/2\Delta t}$
- $P_\varepsilon(\hat{\tau} \geq \Delta t) = O(\varepsilon)$
- $E_\varepsilon[(V(\hat{X}_{\hat{\tau}}, \hat{\tau}) - V(x, 0)) 1_{\hat{\tau} < \Delta t}] = O(\varepsilon)$

Table: Stochastic optimization parameters

Δt	0.01	Time step .
T	1.0	Final time.
σ	0.3	SDE coefficient
r	0.07	SDE coefficient
X_0	100	Initial position for the SDE
K	100	Strike price in the payoff $V(x, t) = e^{-r t} (K - x)_+$
S_{min}	60	Lower bound for the parameter S
S_{max}	90	Upper bound for the parameter S
λ_{min}	-0.01	Lower bound for the parameter λ
λ_{max}	0.7	Upper bound for the parameter λ
L	50	The number of paths started at position X_0 at time 0
S_{path}	50	The number of samples of \bar{D} per path hitting the boundary

Table: Initial condition for the parameters (S, λ)

S_0	80.7642	Initial condition for the parameters.
S_1	70.8288	Initial condition for the parameters.
S_2	79.0376	Initial condition for the parameters.
S_3	76.8769	Initial condition for the parameters.
λ_0	0.556173	Initial condition for the parameters.
λ_1	0.279206	Initial condition for the parameters.
λ_2	0.0235858	Initial condition for the parameters.
λ_3	0.286322	Initial condition for the parameters.

Multidimensional simulation, $d > 1$.

Assume $x^P = 0$ y $\bar{a} = 0$. **Linear change of coordinates:**

- $\bar{X}_s = AZ_s$
- $Z_s - Z_0$ is a Brownian motion in \mathbb{R}^d
- $\bar{h}(\bar{X}_s, s) = 0$ if and only if $e_1^T Z_s = \beta s$.

The stopping time only depends on the new first coordinate and the stopping time is independent of the remaining coordinates. Requirements:

- $AA^T = \bar{\sigma} \bar{\sigma}^T$, **orthogonal columns with the inner product $(\bar{\sigma} \bar{\sigma}^T)^{-1}$**
- $Ae_1 = \frac{1}{\sqrt{\nabla h^T \bar{\sigma} \bar{\sigma}^T \nabla h}} \bar{\sigma} \bar{\sigma}^T \nabla h$, **Conormal vector.**
- $\beta = -\frac{h_t}{\sqrt{\nabla h^T \bar{\sigma} \bar{\sigma}^T \nabla h}}$.

The stopping time for $e_1^T Z_s$ is denoted by $\bar{\tau}$. Define

- $e_i^T Z_{\bar{\tau}} = e_i^T Z_0 + \sqrt{\bar{\tau}} N_i$, $i \neq 1$
- $(\bar{X}_{\bar{\tau}}, \bar{\tau}) = (AZ_{\bar{\tau}}, \bar{\tau})$

with N_i standard normal variables.