Monte Carlo simulation of Heston model

Aurélien Perez & Tina Truong, 2025

Heston model

$$dS_t = rS_t dt + \sqrt{v_t} S_t d\hat{Z}_t \tag{1}$$

$$dv_t = \kappa(\theta - v_t)dt + \sigma\sqrt{v_t}dW_t \tag{2}$$

$$\hat{Z}_t = \rho W_t + \sqrt{1 - \rho^2} Z_t \tag{3}$$

where:

- the spot values $S_0 = 1$ and $v_0 = 0.1$,
- r is the risk-free interest rate, we assume r = 0,
- κ is the mean reversion rate of the volatility,
- θ is the long-term volatility,
- σ is the volatility of volatility,
- W_t and Z_t are independent Brownian motions

Contents

- I. GPU methods
- II. Euler
- III. Exact
- IV. Almost Exact
- V. Results comparisons

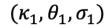
I. GPU methods



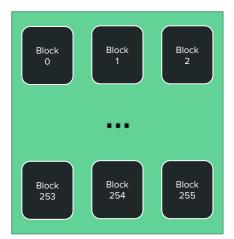
For each parameter set κ , θ , σ , we simulate 2^18 trajectories 1024 Threads per block

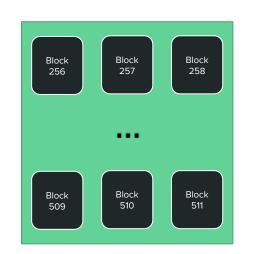
- Simulates 1 trajectory (Euler, Exact or Almost Exact scheme)
- Computes discounted payoff
- Stores result in shared memory (R1s, R2s)

$$(\kappa_0, \theta_0, \sigma_0)$$









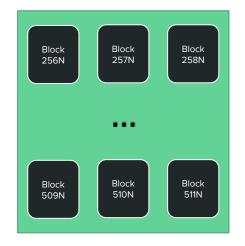
Shared memory reduction per block (sum of 1024 payoffs)

Thread 0 of block → atomicAdd to:

- d_sum[triplet_id]
- d sum2[triplet id]

Final payoff mean and variance (computed outside the kernel)

$$(\kappa_N, \theta_N, \sigma_N)$$

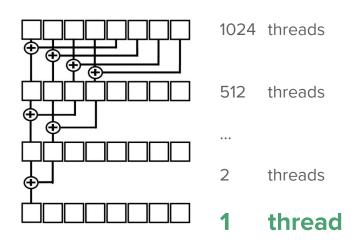


atomicAdd to d_sum[1], d_sum2[1]

Block-Level Reduction



```
// reduction
__syncthreads();
int i = blockDim.x / 2;
while (i!=0){
    if (threadIdx.x < i){</pre>
        R1s[threadIdx.x] += R1s[threadIdx.x + i];
        R2s[threadIdx.x] += R2s[threadIdx.x + i];
    __syncthreads();
    i /= 2;
if (threadIdx.x == 0){
    atomicAdd(&d_sum[triplet_idx], R1s[0]);
    atomicAdd(&d_sum2[triplet_idx], R2s[0]);
```



Atomic add to pass from GPU to CPU and avoid collision.

II. Euler method

I. Euler method

$$\begin{split} S_{t+\Delta t} &= S_t + r S_t \Delta t + \sqrt{v_t} \, S_t \, \sqrt{\Delta t} \left(\rho G_1 + \sqrt{1 - \rho^2} G_2 \right) \\ v_{t+\Delta t} &= g \Big(v_t + \kappa (\theta - v_t) \Delta t + \sigma \, \sqrt{v_t \Delta t} G_1 \Big) \end{split}$$

```
// Euler Scheme kernel
qlobal void MC Heston Euler kernel(float S 0, float v 0, float r, float rho, float sqrt dt,
     float K, int N, curandState *state, float *d sum, float *d sum2, int n triplets, HestonParam *d params) {
    int idx = blockIdx.x * blockDim.x + threadIdx.x;
    int total threads = n triplets * N TRAJ;
    if (idx >= total_threads) return;
    int triplet idx = idx >> POW TRAJ;
    if (triplet idx >= n triplets) return;
    shared HestonParam p shared;
    if (threadIdx.x == 0) p_shared = d_params[triplet_idx];
    __syncthreads();
    HestonParam p = p_shared;
    float kappa = p.kappa, theta = p.theta, sigma = p.sigma;
    float dt = sqrt_dt * sqrt_dt;
    float S = S_0;
    float v = v_0, v_next;
    curandState localState = state[idx];
    extern shared float A[];
    float* R1s, * R2s;
    R1s = A;
    R2s = R1s + blockDim.x;
    for (int i = 0; i < N; i++){
        float2 G = curand normal2(&localState);
        v \text{ next} = fmaxf(v + kappa*(theta - v)*dt + sigma * sgrtf(v) * sgrt dt * G.x, 0.0f);
        S += r * S * dt + sqrtf(v) * S * sqrt_dt * (rho * G.x + sqrtf(1.0f - rho*rho) * G.y);
        v = v \text{ next};
    R1s[threadIdx.x] = expf(-r * dt * N) * fmaxf(0.0f, S-K)/N_TRAJ ;
    R2s[threadIdx.x] = R1s[threadIdx.x] * R1s[threadIdx.x] * N TRAJ;
```

III. Exact method

II. Exact method

step 1. Within a for loop on time steps, we define

$$d=2\kappa\theta/\sigma^2, \quad \lambda=rac{2\kappa e^{-\kappa\Delta t}v_t}{\sigma^2(1-e^{-\kappa\Delta t})}, \quad N=\mathcal{P}(\lambda), \quad ext{with } \mathcal{P} ext{ simulated by curand_poisson}$$

and denoting $\mathcal{G}(\alpha)$ the standard gamma distribution whose simulation is presented in [6]

$$v_{t+\Delta t} = \frac{\sigma^2 (1 - e^{-\kappa \Delta t})}{2\kappa} \mathcal{G}(d+N).$$

- step 2. The integral $\int_0^1 v_s ds$ is stored in a variable vI set to zero before the for loop on time steps. Then, in the for loop we update vI+=0.5 * $(v_t + v_{t+\Delta t})\Delta t$.
- step 3. Once we finish the for loop, we compute $\int_0^1 \sqrt{v_s} dW_s$, using the expression

$$\int_0^1 \sqrt{v_s} dW_s = rac{1}{\sigma} \left(v_1 - v_0 - \kappa heta + \kappa exttt{vI}
ight)$$

Then we compute

$$m=-0.5$$
vI $+
ho\int_0^1\sqrt{v_s}dW_s,\quad \Sigma^2=(1-
ho^2)$ vI

and we set $S_1 = \exp(m + \Sigma G)$ where G is a standard normal random variable independent from (G_1, G_2) .

```
__device__ float rgamma_sup1(curandState *state, float alpha) {
    float d = alpha - 1.0f / 3.0f;
    float c = 1.0f / sqrtf(9.0f * d);
   float z, u, x, v;
   while (true) {
        z = curand_normal(state);
        u = curand uniform(state);
        x = 1.0f + c * z;
       V = X * X * X;
        if (z > -1.0f / c \&\& logf(u) < (0.5f * z * z + d - d * v + d * logf(v)))
            return d * v;
__device__ float rgamma_inf1(curandState *state, float alpha) {
    float u = curand uniform(state);
    return rgamma_sup1(state, alpha + 1.0f) * powf(u, 1.0f / alpha);
// Gamma distribution sampling (Marsaglia-Tsang)
device float rgamma(curandState *state, float alpha) {
    if (alpha < 1.0f) {</pre>
        return rgamma_inf1(state, alpha);
        return rgamma sup1(state, alpha);
```

```
// Exact scheme kernel (Broadie-Kaya)
__global__ void MC_Heston_Exact_kernel(float S_0, float v_0, float r, float rho, float sqrt_dt,
     float K, int N, curandState *state, float *d_sum, float *d_sum2, int n_triplets, HestonParam *d_params) {
    int idx = blockIdx.x * blockDim.x + threadIdx.x;
    int total_threads = n_triplets * N_TRAJ;
   if (idx >= total_threads) return;
   int triplet idx = idx >> POW TRAJ; // idx / 2^POW TRAJ
                  0 1 2 3 4 ... 2^18 - 1 2^18 2^18+1 ...
   // triplet idx 0 0 0 0 0 ... 0 1 1 ...
    if (triplet_idx >= n_triplets) return;
    shared HestonParam p shared;
    if (threadIdx.x == 0) p shared = d params[triplet idx];
    __syncthreads();
   HestonParam p = p_shared;
    float kappa = p.kappa, theta = p.theta, sigma = p.sigma;
    float sigma2 = sigma * sigma;
   float dt = sqrt_dt * sqrt_dt;
   float d = 2.0f * kappa * theta / sigma2;
   float v = v_0, v_next, vI = 0.0f, S = S_0;
   curandState localState = state[idx];
    for (int i = 0; i < N; i++) {
        float lambda = (2.0f * kappa * expf(-kappa * dt) * v) / (sigma2 * (1.0f - expf(-kappa * dt)));
       int N pois = curand poisson(&localState, lambda);
       float gamma = rgamma(&localState, d + N_pois);
       v_next = (sigma2 * (1.0f - expf(-kappa * dt)) / (2.0f * kappa)) * gamma;
       vI += 0.5f * dt * (v_next + v);
       v = v \text{ next};
    float m = -0.5f * vI + (rho / sigma) * (v - v_0 - kappa * theta + kappa * vI);
    float Sigma = sgrtf((1.0f - rho * rho) * vI);
   float2 G = curand normal2(&localState);
   S = S @ * expf(m + Sigma * G.x);
   extern __shared__ float A[];
   float* R1s, * R2s;
   R1s = A;
   R2s = R1s + blockDim.x;
   R1s[threadIdx.x] = expf(-r * dt * N) * fmaxf(0.0f, S-K)/N_TRAJ ;
   R2s[threadIdx.x] = R1s[threadIdx.x] * R1s[threadIdx.x] * N TRAJ;
```

IV. Almost exact method

III. Almost exact method

$$\log(S_{t+\Delta t}) = \log(S_t) + k_0 + k_1 v_t + k_2 v_{t+\Delta t} + \sqrt{(1-\rho^2)v_t \Delta t} \left(\rho G_1 + \sqrt{1-\rho^2} G_2\right)$$

with

$$k_0 = \left(-\frac{\rho}{\sigma}\kappa\theta\right)\Delta t$$
$$k_1 = \left(\frac{\rho\kappa}{\sigma} - 0.5\right)\Delta t - \frac{\rho}{\sigma}$$
$$k_2 = \frac{\rho}{\sigma}$$

where v_t is simulated as in the exact scheme.

```
// Almost Exact Scheme kernel (Haastrecht-Pelsser)
__global__ void MC_Heston_Almost_Exact_kernel(float S_0, float v_0, float r, float rho, float sqrt_dt,
     float K, int N, curandState *state, float *d_sum, float *d_sum2, int n_triplets, HestonParam *d_params) {
    int idx = blockIdx.x * blockDim.x + threadIdx.x;
    int total threads = n triplets * N TRAJ;
    if (idx >= total_threads) return;
    int triplet_idx = idx >> POW_TRAJ;
    if (triplet_idx >= n_triplets) return;
    __shared__ HestonParam p_shared;
    if (threadIdx.x == 0) p_shared = d_params[triplet_idx];
    __syncthreads();
    HestonParam p = p_shared;
    float kappa = p.kappa, theta = p.theta, sigma = p.sigma;
    float sigma2 = sigma * sigma;
    float dt = sqrt_dt * sqrt_dt;
    float d = 2.0f * kappa * theta / sigma2;
    float S = logf(S_0);
    float v = v_0;
    float v_next;
    float rhosigma = rho / sigma;
    float k0 = (-rhosigma * kappa * theta) * dt;
    float k1 = (rhosigma * kappa - 0.5f) * dt - rhosigma;
    float k2 = rhosigma;
    curandState localState = state[idx];
    for (int i = 0; i < N; i++){
        float2 G = curand normal2(&localState);
        float lambda = (2.0f * kappa * expf(-kappa * dt) * v) / (sigma2 * (1.0f - expf(-kappa * dt)));
        int N_pois = curand_poisson(&localState, lambda);
        float gamma = rgamma(&localState, d + N_pois);
        v_next = (sigma2 * (1.0f - expf(-kappa * dt)) / (2.0f * kappa)) * gamma;
       S += k0 + k1 * v + k2 * v_next + sqrtf((1.0f - rho*rho) * v) * sqrt_dt * (rho * G.x + sqrtf(1.0f - rho*rho) * G.y);
        v = v_next;
   S = \exp f(S);
    extern __shared__ float A[];
    float* R1s, * R2s;
    R1s = A;
   R2s = R1s + blockDim.x:
   R1s[threadIdx.x] = expf(-r * dt * N) * fmaxf(0.0f, S-K)/N_TRAJ;
   R2s[threadIdx.x] = R1s[threadIdx.x] * R1s[threadIdx.x] * N_TRAJ;
```

V. Results comparison

Execution speed / triplet

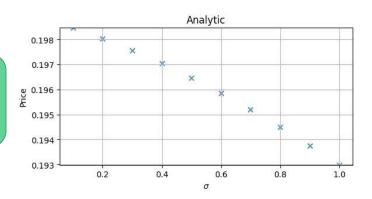
Exact 436.7 ms

Almost Exact 437.01 ms

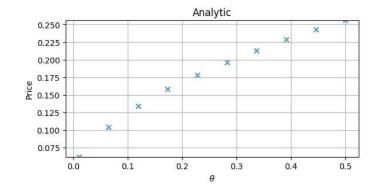
Euler 5.4 ms

Call Price Sensitivity to σ (fixed $\kappa \approx 5.60$, $\theta \approx 0.28$)

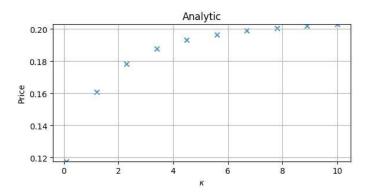
Analytical prices by inverting Fourier transform



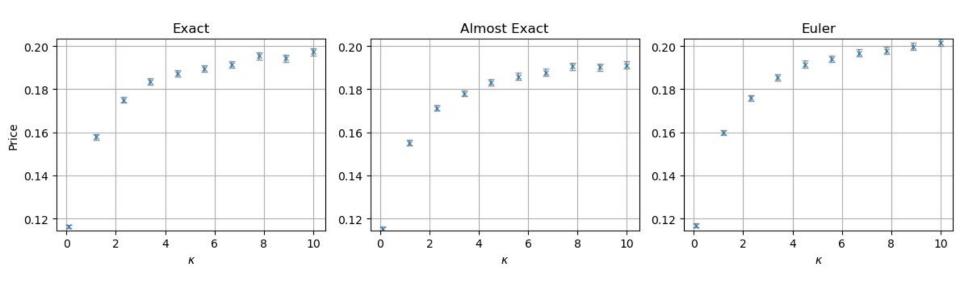
Call Price Sensitivity to θ (fixed $\kappa \approx 5.60$, $\sigma \approx 0.50$)



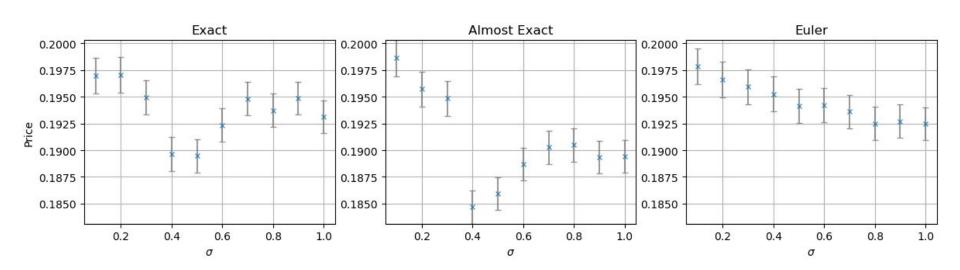
Call Price Sensitivity to κ (fixed $\theta \approx 0.28$, $\sigma \approx 0.50$)



Call Price Sensitivity to κ (fixed $\theta \approx 0.28$, $\sigma \approx 0.50$)



Call Price Sensitivity to σ (fixed $\kappa \approx 5.60$, $\theta \approx 0.28$)



Call Price Sensitivity to θ (fixed $\kappa \approx 5.60$, $\sigma \approx 0.50$)

