Volatility and Correlation Analysis of Financial Market Data

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March 11, 2019

Overview

- Introduction
- ② High-Frequency Prices and Inference
- 3 Open, Close, High, Low Prices
 - Galerkin solution
 - Small-time Analytic Solution and Gap-Fill Approximation

Data type: High-Frequency Prices

Longitudinal view

Time-stamp	Volume	Price	
2011-08-01 03:02:52.112434	1000	179.6400	
2011-08-01 03:02:53.752456	1000	179.6200	
2011-08-01 03:02:53.900010	1000	179.6300	
2011-08-01 03:02:54.103493	1000	179.6050	
2011-08-01 03:02:54.343493	1000	179.6700	

Cross-sectional view: the bid-ask spread

Volume	Price		
101	179.6400		
203	179.6200		
305	179.6100		
500	179.6000		

Data type: Open, Close, High, Low prices





High-frequency returns and difficulties

- As sampling frequency approaches transaction-by-transaction frequency, irregular spacing between transactions and discreteness in transaction prices (such as the bid-ask spread) become dominant features of the data [10].
- These confounding effects in estimating volatility are generally termed microstructure noise, which is
 - independent of sampling interval
 - dominant over short time (< 5 minutes)
 - not cumulative (prices at later times are not affected)

Past work and model-bashed incoherence

 Attempts have been made to fit discrete time state-space models to high-frequency data and estimate price volatility [4, 2]

$$y_{t+\Delta t} = F_t \theta_t + C_t W_{t,1}$$

$$\theta_{t+\Delta t} = G_t \theta_t + V_t W_{t,1}.$$

- As formulated, such types of models are are not robust with respect to the specification of the sampling interval [7, 2, 12].
- Inference over shrinking sampling interval Δt leads to **incoherence**.
- Most recent work has focused on the realized variance [6] estimator:

$$RV_T := \sum_{i=1}^{N/\Delta t} r_{t_i}^2,$$

where $r_{t_i}^2$ is the squared *return* observed over $[t_{i-1}, t_i]$.

Georgi Dinolov (UCSC) March 11, 2019 6 / 66

Realized variance estimator

In the absence of microstructure noise,

$$RV_t = \sum_{i=1}^{N/\Delta t} r_{t_i}^2 \rightarrow \int_0^T \sigma^2(X_t, t) dt$$

- Corrections for the presence of microstructure noise have been proposed in the form of:
 - sampling sparser grids and averaging [11],
 - combining estimators based on sub-sampled data at different frequencies [1],
 - kernel-based estimation [8, 3].
- Disadvantages: averages out information; no inherent model-based projection forward in time.

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A filtering-based approach

modeling microstructure noise

Given the **true** price S_j at sample index j, we model the **measured** price P_j as contaminated by noise due to the bid-ask spread D:

$$P_j = S_j + \nu_j, \qquad \qquad \nu_j \sim U[-D/2, D/2],$$

A first-order Taylor expansion of the **observed log price** $log(P_j)$ produces

$$log(P_j) \approx log(S_j) + \frac{1}{S_j} \nu_j.$$

We further model the noise term with a Gaussian distribution

$$\zeta_j := rac{1}{S_j}
u_j \sim N\left(0, rac{D}{4Q^2}
ight)$$

where Q is an order-magnitude approximation of the true price S_i .

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Stochastic evolution of price and volatility

the idealized joint log-price $\log(\hat{S}_t)$ and log-volatility $\log(\hat{\sigma}_{t,1}), \log(\hat{\sigma}_{t,2})$ diffusion process as the system of SDEs:

$$\begin{split} d\log(\hat{S}_t) &= \hat{\mu} \, dt + \sqrt{\hat{\sigma}_{t,1} \hat{\sigma}_{t,2}} \, \sqrt{dt} \hat{\epsilon}_t + dJ_t, \\ d\log(\hat{\sigma}_{t,1}) &= -\hat{\theta}_1(\log(\hat{\sigma}_{t,1}) - \hat{\alpha}) \, dt + \hat{\tau}_1 \, \sqrt{dt} \hat{\epsilon}_{t,1}, \\ d\log(\hat{\sigma}_{t,2}) &= -\hat{\theta}_2(\log(\hat{\sigma}_{t,2}) - \hat{\alpha}) \, dt + \hat{\tau}_2 \, \sqrt{dt} \hat{\epsilon}_{t,2}. \end{split}$$

- $\log(\hat{\sigma}_{t,1})$ evolves stochastically with long time scale $1/\hat{\theta}_1$,
- $\log(\hat{\sigma}_{t,2})$ evolves stochastically with short time scale $1/\hat{\theta}_2$,
- \bullet dJ_t is a compound Poisson process modeling jumps,
- $\hat{\epsilon}_t, \hat{\epsilon}_{t,1}, \hat{\epsilon}_{t,2}$ are Normally distributed with

$$\mathsf{E}\left[\hat{\epsilon}_t\hat{\epsilon}_{t,2}\right] = \rho, \qquad \quad \mathsf{E}\left[\hat{\epsilon}_t\hat{\epsilon}_{t,2}\right] = 0, \qquad \quad \mathsf{E}\left[\hat{\epsilon}_{t,1}\hat{\epsilon}_{t,2}\right] = 0.$$

Discrete interpretation of the formulation

Defining the log of the observed price P_j as $Y_j := \log(P_j)$

$$Y_{j} = \log(S_{j}) + \zeta_{j},$$

$$\log(S_{j}) = \log(S_{j-1}) + \mu(\Delta) + \sqrt{\sigma_{j,1}\sigma_{j,2}} \,\epsilon_{j} + J_{j}(\Delta),$$

$$\log(\sigma_{j+1,1}) = \alpha(\Delta) + \theta_{1}(\Delta) \left\{ \log(\sigma_{j,1}) - \alpha(\Delta) \right\} + \tau_{1}(\Delta) \,\epsilon_{j,1},$$

$$\log(\sigma_{j+1,2}) = \alpha(\Delta) + \theta_{2}(\Delta) \left\{ \log(\sigma_{j,2}) - \alpha(\Delta) \right\} + \tau_{2}(\Delta) \,\epsilon_{j,2}.$$

$$\sigma_{j+1,i} = \hat{\sigma}_{(j+1)\Delta,i}\sqrt{\Delta}, \quad S_j = \hat{S}_{j\Delta}, \quad J_j(\Delta) = J((j+1)\Delta) - J(j\Delta)$$

$$lpha(\Delta) = \hat{lpha} + rac{1}{2}\log(\Delta), \quad \mu(\Delta) = \hat{\mu}\Delta,$$
 $au_i(\Delta) = \hat{ au}_i\sqrt{rac{1 - \exp\left\{-2\hat{ heta}_i\Delta\right\}}{2\hat{ heta}_i}}, \quad heta_i(\Delta) = \exp\left\{-\hat{ heta}_i\Delta\right\}.$

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Prior formulation

To ensure that the model estimation is coherent across choices of sampling periods $\boldsymbol{\Delta},$ we

- define the first two moments of each continuous-time parameter,
- use the Delta Method to calculate the first two moments of each discrete-time parameter,
- 3 define the prior for each discrete-time parameter.

Computation

• We separate the standard innovation term from volatility factors $\sigma_{j,1}$ and $\sigma_{j,2}$ by considering the transformation

$$\log(S_{j}) = \log(S_{j-1}) + \mu(\Delta) + \sqrt{\sigma_{j,1}\sigma_{j,2}} \, \epsilon_{j} + J_{j}(\Delta)$$

$$\leftrightarrow \underbrace{\log\left[\left|\log(S_{j}/S_{j-1}) - \mu(\Delta) - J_{j}(\Delta)\right|\right]}_{y_{j}^{*}} = \underbrace{\frac{1}{2} \underbrace{\log(\sigma_{j,1})}_{h_{j,1}} + \underbrace{\frac{1}{2} \underbrace{\log(\sigma_{j,2})}_{h_{j,2}}}_{h_{j,2}} + \underbrace{\underbrace{\log(\epsilon_{j}^{2})/2}_{\epsilon_{*}^{*}}.$$

Computation

• We separate the standard innovation term from volatility factors $\sigma_{j,1}$ and $\sigma_{j,2}$ by considering the transformation

$$\log(S_{j}) = \log(S_{j-1}) + \mu(\Delta) + \sqrt{\sigma_{j,1}\sigma_{j,2}} \, \epsilon_{j} + J_{j}(\Delta)$$

$$\leftrightarrow \underbrace{\log\left[\left|\log(S_{j}/S_{j-1}) - \mu(\Delta) - J_{j}(\Delta)\right|\right]}_{y_{j}^{*}} = \frac{1}{2} \underbrace{\log(\sigma_{j,1})}_{h_{j,1}} + \underbrace{\frac{1}{2} \underbrace{\log(\sigma_{j,2})}_{h_{j,2}}}_{e_{j}^{*}}.$$

• We approximate ϵ_i^* as a mixture of Normals

$$\epsilon_j^* = \log(\epsilon_j^2)/2 \sim \sum_{l=1}^{10} p_l N\left(\frac{m_l}{2}, \frac{v_l^2}{4}\right).$$

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Computation, continued

The joint distribution $(\epsilon_j^*, \epsilon_{j,2})$ conditional on the latent mixture element γ_j becomes

$$\begin{split} p(\epsilon_{j}^{*}, \epsilon_{j,2} | \gamma_{j}) &= p(\epsilon_{j,2} | \epsilon_{j}^{*}, \gamma_{j}) p(\epsilon_{j}^{*} | \gamma_{j}) \\ &= p(\epsilon_{j,2} | \underbrace{d_{j} exp(\epsilon_{j}^{*})}_{\epsilon_{j}}, \gamma_{j}) p(\epsilon_{j}^{*} | \gamma_{j}) \\ &= N\left(\epsilon_{j,2} \left| d_{j} exp(\epsilon_{j}^{*}), (1 - \rho^{2}) \right.\right) N\left(\epsilon_{j}^{*} \left| \frac{m_{\gamma_{j}}}{2}, \frac{v_{\gamma_{j}}^{2}}{4} \right.\right), \end{split}$$

where d_j is the sign of ϵ_j .

Computation, continued

The joint distribution $(\epsilon_j^*, \epsilon_{j,2})$ conditional on the latent mixture element γ_j becomes

$$\begin{split} p(\epsilon_{j}^{*}, \epsilon_{j,2} | \gamma_{j}) &= p(\epsilon_{j,2} | \epsilon_{j}^{*}, \gamma_{j}) p(\epsilon_{j}^{*} | \gamma_{j}) \\ &= p(\epsilon_{j,2} | \underbrace{d_{j} exp(\epsilon_{j}^{*})}_{\epsilon_{j}}, \gamma_{j}) p(\epsilon_{j}^{*} | \gamma_{j}) \\ &= N\left(\epsilon_{j,2} \left| d_{j} exp(\epsilon_{j}^{*}), (1 - \rho^{2}) \right.\right) N\left(\epsilon_{j}^{*} \left| \frac{m_{\gamma_{j}}}{2}, \frac{v_{\gamma_{j}}^{2}}{4} \right.\right), \end{split}$$

where d_j is the sign of ϵ_j . The joint distribution becomes a linear combination of independent Normal random variables when we replace the nonlinear $exp(\epsilon_j^*)$ term with the first-order approximation

$$exp(\epsilon_j^*)|\gamma_jpprox exp(m_{\gamma_j}/2)(a_{\gamma_j}+b_{\gamma_j}(2\epsilon_j^*-m_{\gamma_j}))$$

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Blocked Gibbs sampler with predominantly Metropolis-Hastings and Forward Filtering Backward Sampling steps

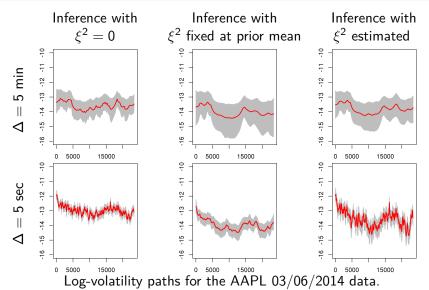
- sample observational parameters (MH)
- 2 sample latent prices (FFBS)
- sample volatility parameters (MH)
- sample latent mixture indicators (Discrete)
- sample volatility paths (FFBS)
- sample jump parameters (MH)
- sample jumps (Conjugate prior)

Estimating integrated volatility

- $\zeta \sim N(0, \xi^2)$
- We consider 300 simulated data sets over 1 trading day
- Coverage : percent of 95% probability/confidence intervals covering the true data-generating integrated volatility $\int \hat{\sigma}_t^2 dt$:

	Sampling period			
	60 sec	30 sec	15 sec	5 sec
Inference with $\xi^2 = 0$	72	28	3	0
Inference with ξ^2 fixed at prior mean	79	57	23	0
Inference with ξ^2 estimated	91	92	96	97
Inference with kernel-based estimator	51	48	59	76

Results for real data example



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 16 / 66

Bivariate Open, Close, High, Low Prices

- Work has been done in the univariate case
- No equivalent results in the bivariate setting
- There exists only a single estimator in the literature which used bivariate OCHL to estimate asset correlation [9] (Rogers estimator)

Model formulation

We consider a two-dimensional correlated Brownian motion:

$$X(t) = x_0 + \mu_x t + \sigma_x W_x(t)$$

$$Y(t) = y_0 + \mu_y t + \sigma_y W_y(t)$$

where $W_x(t)$ and $W_y(t)$ are correlated standard Brownian motions with $\text{Cov}(W_x(t),W_x(t))=\rho t$.

• We seek the 6-dimensional joint probability density function for the pair (X(t), Y(t)) and the random variables $M_X(t) = \max_{0 \le s \le t} X(s)$, $m_X(t) = \min_{0 \le s \le t} X(s)$, $M_Y(t) = \max_{0 \le s \le t} Y(s)$, $m_Y(t) = \min_{0 \le s \le t} Y(s)$.

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Transition density

To calculate the joint density

$$p(X(t) = x, Y(t) = y,$$

 $m_X(t) = a_X, M_X(t) = b_X, m_Y(t) = a_Y, M_Y(t) = b_Y),$

we first study cumulative-like distribution

$$p(X(t) = x, Y(t) = y,$$

 $m_X(t) \ge a_X, M_X(t) \le b_X, m_Y(t) \ge a_Y, M_Y(t) \le b_Y,$

which is governed by a Fokker-Planck equation with absorbing boundaries:

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Governing Fokker-Planck equation

Abbreviating

$$q(x, y, t) := p(X(t) = x, Y(t) = y,$$

 $m_X(t) \ge a_X, M_X(t) \le b_X, m_Y(t) \ge a_Y, M_Y(t) \le b_Y,$

the FP equation is

$$\begin{split} \frac{\partial}{\partial t'} q(x, y, t') &= -\mu_x \frac{\partial}{\partial x} q(x, y, t') - \mu_y \frac{\partial}{\partial y} q(x, y, t') + \\ &\frac{1}{2} \sigma_x^2 \frac{\partial^2}{\partial x^2} q(x, y, t') + \rho \sigma_x \sigma_y \frac{\partial^2}{\partial x \partial y} q(x, y, t') + \frac{1}{2} \sigma_y^2 \frac{\partial^2}{\partial y^2} q(x, y, t'), \end{split}$$

$$q(a_x, y, t') = q(b_x, y, t') = q(x, a_y, t') = q(x, b_y, t') = 0, \quad 0 < t' \le t,$$

$$q(x, y, 0) = \delta(x_0 - x)\delta(y_0 - y)$$

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Normalized problem

We can introduce a series of transformations to solve the equivalent initial-boundary problem

$$\begin{split} \frac{\partial}{\partial \tilde{t}} q(\tilde{x}, \tilde{y}, \tilde{t}) &= \tilde{\mathcal{L}} q(\tilde{x}, \tilde{y}, \tilde{y}), \\ \tilde{\mathcal{L}} &= \frac{1}{2} \frac{\partial^2}{\partial \tilde{x}^2} + \rho \sigma_{\tilde{y}} \frac{\partial^2}{\partial \tilde{x} \partial \tilde{y}} + \frac{1}{2} \sigma_{\tilde{y}}^2 \frac{\partial^2}{\partial \tilde{y}^2}, \quad \tilde{\Omega} := (0, 1) \times (0, 1) \end{split}$$

where $0 < \sigma_{\tilde{v}}^2 \le 1$.

 This new parameterization makes the computational domain invariant to data/parameter combinations.

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Joint density is given by the fourth derivative with respect to boundaries

Denoting the joint density of the diffusion process and the attained extrema over the interval t as $f(x, y, a_x, b_x, a_y, b_y)$,

$$\frac{\partial^4}{\partial a_x \partial b_x \partial a_y \partial b_y} q(x, y, t) = f(x, y, a_x, b_x, a_y, b_y).$$

• The finite difference approximation of $\partial^4/\partial a_x \partial b_x \partial a_y \partial b_y$ introduces a fundamental limitation due to finite precision round-off errors when the analytic solution q(x,y,t) is not available

$$\mathcal{O}\left(rac{arepsilon_{\mathit{mach}}}{arepsilon^4}
ight) o \infty \ \ \ \mathsf{as} \ \ arepsilon o 0,$$

where ε is the step size of the numerical differentiation with respect to the boundaries.

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Approach 1: Finite Difference approximation to FP equation

Requires the solution to a system of ODEs:

$$\dot{c}(\tilde{t}) = Bc(\tilde{t})$$
 $\Rightarrow c(\tilde{t}) = \exp(B\tilde{t}) c(0),$

where the sparse system matrix B can be computed once and stored for a regular grid over $\tilde{\Omega}$ with finite step size h.

$$B = \frac{1}{2} \frac{1}{h^2} B_{\tilde{x}, \tilde{x}} + \rho \sigma_{\tilde{y}} \frac{1}{4h^2} B_{\tilde{x}, \tilde{y}} + \frac{1}{2} \sigma_{\tilde{y}}^2 \frac{1}{h^2} B_{\tilde{y}, \tilde{y}}.$$

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Approach 1: Finite Difference approximation to FP equation

Letting k := 1/h and b denote the boundary parameters,

FD approximation is fast but also limited by irregular truncation errors

$$q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t}|b) - q(\tilde{x}, \tilde{y}, \tilde{t}|b) = \underbrace{\left(\frac{1}{k}\right)^{\alpha} \mathsf{F}_{\mathsf{reg}}(b)}_{\mathsf{smooth in } b, \ \alpha > 0}$$

+
$$\underbrace{\left(\frac{1}{k}\right)^{\beta}\mathsf{F}_{irreg}(b)}_{\text{continuous but not differentiable w.r.t }b,\,\beta>0} + \underbrace{\varepsilon_{mach}\mathsf{F}_{round}(b)}_{\text{behaves as a R.V}}$$

 With linear interpolation which arises when function arguments are not on grid points with respect to boundary differentiation

$$\mathcal{O}(\mathsf{F}_{\mathsf{irreg}}(b)) = \mathsf{k}^{\beta}/\varepsilon \qquad \Rightarrow \left(\frac{1}{\mathsf{k}}\right)^{\beta} \mathsf{F}_{\mathsf{irreg}}(b) o \infty \ \ \mathsf{as} \ \ arepsilon o 0.$$

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Approach 2: Trigonometric expansion of the differential operator

$$q(\tilde{x}, \tilde{y}, \tilde{t}) \approx \sum_{\nu} h_{\nu} \phi_{\nu}(\tilde{x}, \tilde{y}) e^{-\lambda_{\nu} \tilde{t}}.$$

• When $\rho = 0$,

$$q(\tilde{x}, \tilde{y}, \tilde{t}) \approx \sum_{l=1}^{L} \sum_{m=1}^{M} c_{l,m} \sin(2\pi I \tilde{x}) \sin(2\pi m \tilde{y})$$

- the system matrix is diagonal
- no new error is introduced in the time evolution
- truncated solution converges fast

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Approach 2: Trigonometric expansion of the differential operator

• When $\rho \neq 0$, mixing term in the FP equation produces $\cos(2\pi I\tilde{x})\cos(2\pi k\tilde{y})$ terms in the above expansion such that

$$\phi_{\nu}(\tilde{x}, \tilde{y}) = \sum_{l=1}^{L} \sum_{m=1}^{M} c_{l,m,\nu} \sin(2\pi l \, \tilde{x}) \sin(2\pi m \, \tilde{y}) := \Psi(\tilde{x}, \tilde{y})^{T} c_{\nu},$$

- system matrix is dense and convergence is slow,
- new error is introduced in the time evolution

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The Galerkin approach: weak solution to the PDE

We propose a solution $q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t})$ of similar form

$$q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t}) = \sum_{i=0}^{k} c_i(\tilde{t}) \psi_i(\tilde{x}, \tilde{y}),$$

where the basis functions $\psi_i(\tilde{x}, \tilde{y})$ satisfy the boundary conditions on $\tilde{\Omega}$ and they need not be eigenfunctions but capture some essential features of the solution. We also require that all first- and second-order derivatives of $\psi_i(\tilde{x}, \tilde{y})$ are in $L_2(\tilde{\Omega})$.

$$\frac{\partial}{\partial \tilde{t}} q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t}) - \tilde{\mathcal{L}} q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t}) := R_{e}(k),$$
$$q(\tilde{x}, \tilde{y}, 0) - q^{(k)}(\tilde{x}, \tilde{y}, 0) := R_{0}(k).$$

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The Galerkin approach: weak solution to the PDE

The *orthogonality* condition of the Galerkin procedure:

$$\int_{\Omega} R_e(k) \psi_i(\tilde{x}, \tilde{y}) d\tilde{x} d\tilde{y} = 0, \quad \int_{\Omega} R_0(k) \psi_i(\tilde{x}, \tilde{y}) d\tilde{x} d\tilde{y} = 0, \quad i = 0, \dots k,$$

which is equivalent to the weak formulation of the heat problem

$$\left\langle \partial_t q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t}), \psi_i \right\rangle = \left\langle \tilde{\mathcal{L}} q^{(k)}(\tilde{x}, \tilde{y}, \tilde{t}), \psi_i \right\rangle,$$
$$\left\langle q^{(k)}(\tilde{x}, \tilde{y}, 0), \psi_i \right\rangle = \left\langle q(\tilde{x}, \tilde{y}, 0), \psi_i \right\rangle,$$

where $<\cdot,\cdot>$ is the usual inner product in $L_2(\tilde{\Omega})$.

Basis element choice

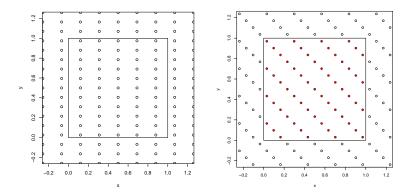
$$\psi_{i}(\tilde{x}, \tilde{y}) = \frac{1}{2\pi\tilde{\sigma}^{2}\sqrt{1-\tilde{\rho}^{2}}}$$

$$\times \exp\left\{-\frac{((\tilde{x}-\tilde{x}_{i})^{2}-2\tilde{\rho}(\tilde{x}-\tilde{x}_{i})(\tilde{y}-\tilde{y}_{i})+(\tilde{y}-\tilde{y}_{i})^{2})}{2(1-\tilde{\rho}^{2})\tilde{\sigma}^{2}}\right\}$$

$$\times \tilde{x}(1-\tilde{x}) \tilde{y}(1-\tilde{y})$$

for some parameters $(\tilde{\rho}, \tilde{\sigma})$ and a collection of nodes $\{(\tilde{x}_i, \tilde{y}_i)\}_{i=0}^k$ which form a grid over $\tilde{\Omega}$.

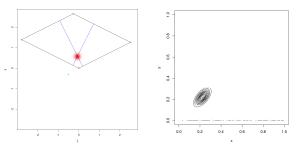
- Align basis elements along the principal axis of the differential operator,
- \bullet Spacing between nodes is I times the bandwidth $\tilde{\sigma}$ in each principal direction
- This scheme is aimed at better resolving the correlation in the solution



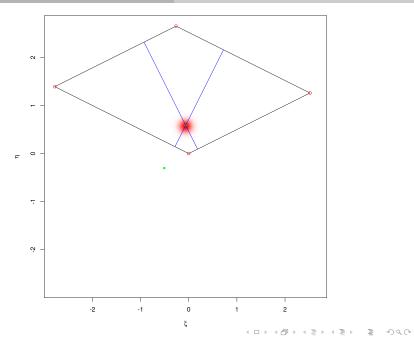
- ullet A sample grid design for I=1, $ilde{\sigma}=0.3$ and $ilde{
 ho}=0.6$.
- **Left:** Grid along coordinates of observed prices within the computational domain (solid black square).
- **Right:** Grid along the major axes of the basis element kernels. The set of final node points $\{(x_i, y_i)\}_{i=0}^k$ is contained within the computational domain and is denoted by the red solid points.

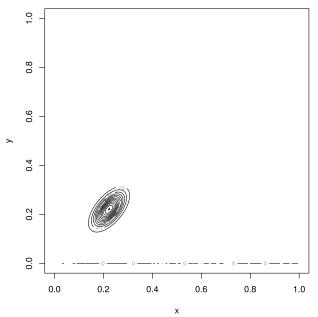
Small-time solution to the Fokker-Planck equation

- A semi-analytic solution for a small time \tilde{t}_{ϵ} allows us to project a smooth function onto our basis family instead of a δ -function
- This reduces the numerical error of the Galerkin solver



- **Left:** Computational domain under the transformation where diffusion parameters are both unity and correlation is zero.
- Right: Contour of small-time solution in original coordinate system.





Consistency Results

Denoting the approximate density due to the Galerkin solution as

$$f^{(k)}(x, y, a_x, b_x, a_y, b_y) = \frac{\partial^4 q^{(k)}(x, y, t)}{\partial a_x \partial b_x \partial a_y \partial b_y},$$

we show that the two integrals below converge in $L_2(\Omega)$:

Lemma (1)

$$\lim_{k \to \infty} \int_{a_x}^{b_x} \int_{a_y}^{b_y} \left(f^{(k)}(x, y, a_x, b_x, a_y, b_y) - f(x, y, a_x, b_x, a_y, b_y) \right)^2 dx dy$$

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Denoting the random variable

$$Z := (X(t), Y(t), m_X(t), M_X(t), m_Y(t), M_Y(t)),$$

we consider the distribution function $F(Z \le z)$ generated by the true solution to the Fokker-Planck equation as well as the distribution $F^{(k)}(Z \le z)$ generated by the Galerkin solution.

Lemma (Convergence in distribution)

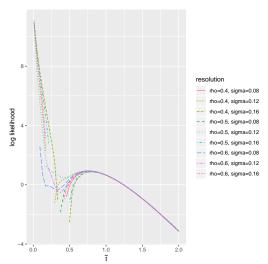
For any
$$z \in Z$$
, $\lim_{k \to \infty} F^{(k)}(Z \le z) = F(Z \le z)$.

Assuming that the maximum likelihood estimate (MLE) (under $F^{(k)}$ for sufficiently large k) is continuous with respect to the data, we can use the convergence result above and Chebyshev's inequality to show that the MLE under $F^{(k)}$ is consistent as the number of basis elements k and number of data points n go to infinity.

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Results and solution behavior for small $ilde{t}$

Here we consider the parameters defining the basis family $(\tilde{
ho}, \tilde{\sigma})$



Simulation Study

We consider 50 simulated data sets with

$$\sigma_x = 1,$$
 $\sigma_y = 1,$ $\mu_x = 0,$ $\mu_y = 0$

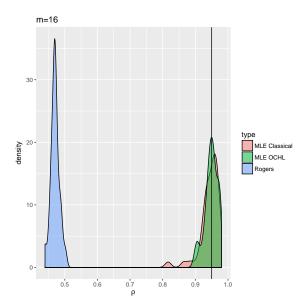
with increasing sample size over several values of ρ . For each case, we consider the ratio of mean-square error of the Galerkin solver compared to that generated by Gaussian likelihoods ignoring the boundaries:

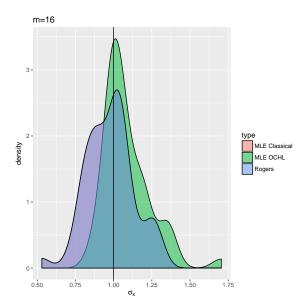
	ho = 0.95			ho=0.60		
	m = 4	m = 8	m = 16	m=4	m = 8	m = 16
$\hat{\sigma}_{x}$	0.475	1.238	1.304	0.203	0.127	0.232
$\hat{\sigma}_y$	0.593	1.040	1.088	0.111	0.120	0.260
$\hat{\rho}$	0.287	0.910	0.445	0.315	0.283	0.463

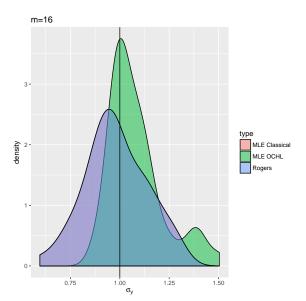
$$ho = 0.0$$
 $m = 4$ $m = 8$ $m = 16$
 ho
 ho

Simulation study

- For data generated with $\rho=0.95$, kernel-density approximations of the repeated-sampling densities of the MLEs are shown. Samples are obtained from the Galerkin likelihood (green) and the classical Gaussian likelihood (red) The data-generating parameters are denoted with the vertical solid line.
- The Rogers estimator (blue) is the only existing correlation estimator for bivariate OCHL data.

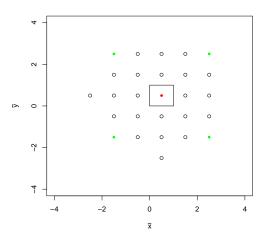






Open, Close, High, Low Prices: Small Time Solution and Gap-Fill Solution

Method of images and analytic differentiability



Symmetry constraint in constructing approximate solutions

A condition weaker than uniqueness which nonetheless restricts the solution space is the symmetry obeyed by the problem. We consider the transformation

$$x^{new} = (a_x + b_x) - x^{old},$$

 $y^{new} = (a_y + b_y) - y^{old}.$

Performing the set of reflections

$$\{2,4,1,3\} \cup \{2,4,3,1\} \cup \{4,2,1,3\} \cup \{4,2,3,1\}$$
.

produces a sum of images where only four elements are differentiable with respect to all four boundaries:

$$\frac{\partial^4 p_{\epsilon}(\tilde{x},\tilde{y},\tilde{t})}{\partial a_X \partial b_X \partial a_y \partial b_y} = \sum_{j'=1}^4 \frac{\partial^4 G(\tilde{x},\tilde{y},\tilde{t}|\tilde{x}_{(j')},\tilde{y}_{(j')})}{\partial a_X \partial b_X \partial a_y \partial b_y}.$$

Calculation of the joint density

We can express the derivatives:

$$\frac{\partial^{4}}{\partial a_{x} \partial b_{x} \partial a_{y} \partial b_{y}} G(\tilde{x}, \tilde{y}, \tilde{t} | \tilde{x}_{(j)}, \tilde{y}_{(j)}) =$$

$$G \cdot \mathcal{C}^{4} \cdot \left(\partial_{a_{x}} \partial_{b_{x}} \partial_{a_{y}} \partial_{b_{y}}\right) \mathcal{P} + G \cdot \mathcal{C}^{3} \cdot \left(\partial_{a_{x}}^{2} b_{x} \partial_{a_{y}} \partial_{b_{y}} + \partial_{a_{x}}^{2} a_{y} \partial_{b_{x}} \partial_{b_{y}} + \partial_{a_{x}}^{2} b_{y} \partial_{a_{x}} \partial_{a_{y}} \partial_{a_{y}} \partial_{b_{x}} \partial_{a_{y}} + \partial_{b_{x}}^{2} b_{y} \partial_{a_{x}} \partial_{a_{y}} \partial_{a_{y}} \partial_{a_{y}} \partial_{a_{y}} \partial_{b_{x}} \partial_{b_{x}}\right) \mathcal{P} + G \cdot \mathcal{C}^{2} \cdot \left(\partial_{a_{x}}^{3} b_{x} a_{y} \partial_{b_{y}} + \partial_{a_{x}}^{2} b_{x} \partial_{a_{y}}^{2} b_{y} \partial_{a_{y}} + \partial_{a_{x}}^{3} b_{x} b_{y} \partial_{a_{y}} + \partial_{a_{x}}^{3} b_{x} b_{y} \partial_{a_{x}} + \partial_{b_{x}}^{2} a_{y} \partial_{b_{y}} \partial_{a_{x}} \partial_{b_{y}}\right) \mathcal{P} + G \cdot \mathcal{C} \cdot \partial_{a_{x}}^{4} b_{x} a_{y} b_{y} \partial_{b_{x}} + \mathcal{P}$$

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Analytic Derivative and higher order terms

Thinking of \tilde{t} as variable allows us to further simplify the expression. Since $\mathcal{C} = \mathcal{O}(1/\tilde{t})$, all three terms $G \cdot \mathcal{C}^3$, $G \cdot \mathcal{C}^2$, and $G \cdot \mathcal{C}$ are $o\left(G \cdot \mathcal{C}^4 \cdot \left(\partial_{a_x} \partial_{b_x} \partial_{a_y} \partial_{b_y}\right) \mathcal{P}\right)$, so that the $G \cdot \mathcal{C}^4$ order term in the derivative dominates the others for sufficiently small \tilde{t} .

$$\frac{\partial^4 p_{\epsilon}(\tilde{x}, \tilde{y}, \tilde{t})}{\partial a_{x} \partial b_{x} \partial a_{y} \partial b_{y}} \approx \sum_{j'=1}^{4} G(\tilde{x}, \tilde{y}, \tilde{t} | \tilde{x}_{(j')}, \tilde{y}_{(j')}) \cdot \mathcal{C}^4 \cdot \left(\partial_{a_{x}} \partial_{b_{x}} \partial_{a_{y}} \partial_{b_{y}}\right) \mathcal{P}_{j'}.$$

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Basis Expansion of Galerkin likelihood: RHS

The likelihood computed with the Galerkin solution as a function of \tilde{t} is of the form

$$\frac{\partial^{4}q_{\mathsf{Galerkin}}(\tilde{x},\tilde{y},\tilde{t})}{\partial a_{x}\partial b_{x}\partial a_{y}\partial b_{y}} = \sum_{k=1}^{K} e^{-\lambda_{k}\tilde{t}} p_{k}^{(4)}(\tilde{t}),$$

where $p_i^{(4)}(\tilde{t})$ is a fourth-order polynomial. This proceeds from the Galerkin solution being dependent on \tilde{t} only through the exponential term: the eigenfunctions of the solution are by design solely functions of (a_x,b_x,a_y,b_y) and (\tilde{x},\tilde{y})

 This suggests a leading-order approximation that is fitted with the Galerkin solver via least squares:

$$f_{\mathsf{Galerkin}}(\tilde{t}) = \tilde{t}^4 \left(\omega_1 e^{-\lambda_1 \tilde{t}} + \omega_2 e^{-\lambda_2 \tilde{t}} \right),$$

$$\log f_{\mathsf{Galerkin}}(\tilde{t}) = \log(\omega_1) + 4 \log(\tilde{t}) - \lambda_1 \tilde{t} + \log\left(1 + \omega_2/\omega_1 e^{-(\lambda_2 - \lambda_1)\tilde{t}}\right)$$

Leading order for small-time solution: LHS

The summand in the small-time likelihood with the greatest $\beta_{j'}$ contributes the most to the truncated small-time solution in the $\tilde{t} \leq 1$ region where the matched solution will be applied. Indexing j' such that $\beta_1 \geq \beta_2 \geq \beta_3 \geq \beta_4$, the small-time log-likelihood is

$$\begin{split} \log \left(\frac{\partial^4 p_{\epsilon}(\tilde{x}, \tilde{y}, \tilde{t})}{\partial a_{x} \partial b_{x} \partial a_{y} \partial b_{y}} \right) &\approx \log(K) - 4.5 \log(\tilde{t}) + \log(c_{1}) - \frac{\beta_{1}}{\tilde{t}} \\ &+ \log \left(1 + \sum_{j \neq 1} \frac{c_{j}}{c_{1}} \exp\left(-\frac{(\beta_{j} - \beta_{1})}{\tilde{t}} \right) \right) \\ &\approx \log(K) - 4.5 \log(\tilde{t}) + \log(c_{1}) - \frac{\beta_{1}}{\tilde{\tau}} + \log\left(1 + \epsilon(\tilde{t}) \right) \end{split}$$

• The proposed matched solution is of the form

$$\log f_{\sf gap}(ilde{t}) = \log(\omega(ilde{t})) - \gamma(ilde{t})\log(ilde{t}) - rac{eta(t)}{ ilde{t}}$$

Georgi Dinolov (UCSC) March 11, 2019 48 / 66

Gap-fill solution

$$\begin{split} \log f_{\mathsf{Galerkin}}(\tilde{t}) &\approx \log(\omega_1) + 4\log(\tilde{t}) - \lambda_1 \tilde{t} + \log\left(1 + \omega_2/\omega_1 e^{-(\lambda_2 - \lambda_1)\tilde{t}}\right) \\ \log f_{\mathsf{small-time}}(\tilde{t}) &\approx \log(K) - 4.5\log(\tilde{t}) - \frac{\beta_1}{\tilde{t}} + \log(c_1) + \log\left(1 + \epsilon(\tilde{t})\right), \\ \log f_{\mathsf{gap}}(\tilde{t}) &= \log(\omega(\tilde{t})) - \gamma(\tilde{t})\log(\tilde{t}) - \frac{\beta(\tilde{t})}{\tilde{t}} \end{split}$$

• At \tilde{t}^* , the left-hand side of the matching condition, the values for these parameters are defined such that they match the small-time solution

$$\omega(\tilde{t}^*) = K, \qquad \gamma(\tilde{t}^*) = 4.5, \qquad \beta(\tilde{t}^*) = \beta_1.$$

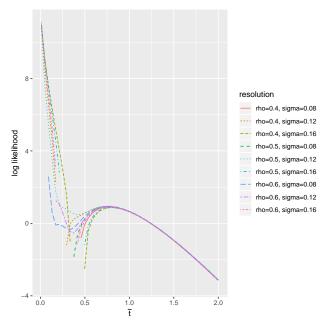
• At \tilde{t}_m , the right-hand side of the matching condition and the maximum of the LS solution, $\omega(\tilde{t}), \gamma(\tilde{t})$, and $\beta(\tilde{t})$ are chosen to match the value, first, and second derivatives of the logarithmic form of the Galerkin approximate log-likelihood.

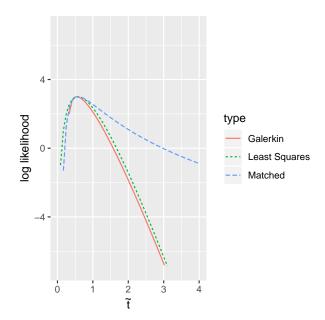
$$\omega(\tilde{t}) = \omega(\tilde{t}^*)e^{-k(\tilde{t}-\tilde{t}^*)} + \omega(\tilde{t}_m) \left(1 - e^{-k(\tilde{t}-\tilde{t}^*)}\right),$$

$$\gamma(\tilde{t}) = \gamma(\tilde{t}^*)e^{-k(\tilde{t}-\tilde{t}^*)} + \gamma(\tilde{t}_m) \left(1 - e^{-k(\tilde{t}-\tilde{t}^*)}\right),$$

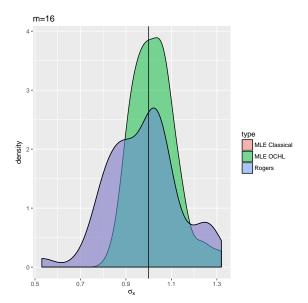
$$\beta(\tilde{t}) = \beta(\tilde{t}^*)e^{-k(\tilde{t}-\tilde{t}^*)} + \beta(\tilde{t}_m) \left(1 - e^{-k(\tilde{t}-\tilde{t}^*)}\right).$$

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Repeated Simulation Results



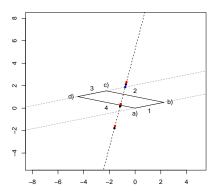
- Multivariate High-Frequency models
- Gaussian Process Emulator for interpolating the OCHL likelihood
- Particle Filter

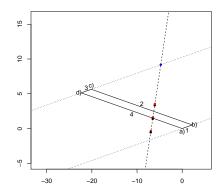


Thank you



Illustration of proof for existence







Prior elicitation example I

Prior for $\theta_i(\Delta)$: The discrete-time autocorrelation coefficient $\theta(\Delta)$ of the volatility process is bounded above by 1 and below by 0 such that $\log(\sigma_j)$ is bounded as $j \to \infty$. Hence, we employ a truncated normal prior for $\theta(\Delta)$,

$$\textit{p}(\theta(\Delta)) \propto \textit{N}\left(\textit{a}_{\theta}(\Delta),\textit{b}_{\theta}^{2}(\Delta)\right)\mathbb{1}_{\left(\theta(\Delta) \in [0,1]\right)},$$



Prior elicitation example II

which leads again to a tractable computational algorithm. Note that because of the truncation,

$$E\left[\theta(\Delta)\right] = a_{\theta}(\Delta) + \frac{\phi\left(-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right) - \phi\left(\frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right)}{\Phi\left(\frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right) - \Phi\left(-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right)} b_{\theta}(\Delta) \tag{1}$$

$$Var\left[\theta(\Delta)\right] = b_{\theta}^{2}(\Delta) \left[1 + \frac{-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\phi\left(-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right) - \frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\phi\left(\frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right)}{\Phi\left(\frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right) - \Phi\left(-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right)} + \left\{\frac{\phi\left(-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right) - \phi\left(\frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right)}{\Phi\left(\frac{1 - a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right) - \Phi\left(-\frac{a_{\theta}(\Delta)}{b_{\theta}(\Delta)}\right)}\right\}^{2} \right] \tag{2}$$

where $\phi(\cdot)$ and $\Phi(\cdot)$ denote the density and the cumulative distribution functions of the standard normal distribution. Now, given the prior mean

Georgi Dinolov (UCSC) March 11, 2019 59 / 66

Prior elicitation example III

 $\hat{a}_{\hat{\theta}}$ and variance $\hat{b}_{\hat{\theta}}^2$ for $\hat{\theta}$, we choose the values of $a_{\theta}(\Delta)$ and $b_{\theta}(\Delta)$ so that the mean and variance of $\theta(\Delta)$ above are approximately equal to the mean and variance of $\exp\left\{-\hat{\theta}\Delta\right\}$. To simplify calculation of the moments of $\exp\left\{-\hat{\theta}\Delta\right\}$ we use a second-order Taylor expansion of $\exp\left\{-\hat{\theta}\Delta\right\}$ to approximate the first two moments of $\theta(\Delta)$ in terms of $\hat{a}_{\hat{\theta}}$ and \hat{b}_{θ}^2 , an approach known as the Delta-Method (e.g., see [5]):

$$\mathsf{E}\left[\exp\left\{-\hat{\theta}\Delta\right\}\right] \approx \exp\left(-\hat{a}_{\hat{\theta}}\Delta\right) \left(1 + \frac{1}{2}\hat{b}_{\hat{\theta}}^2\Delta^2\right),\tag{3}$$

$$\mathsf{E}\left[\exp\left\{-2\hat{\theta}\Delta\right\}\right] \approx \exp\left(-2\hat{a}_{\hat{\theta}}\Delta\right) \left(1 + 2\hat{b}_{\hat{\theta}}^2\Delta^2\right). \tag{4}$$

Using (1), (2), (3), and (4), and by setting

$$\mathsf{E}\left[\theta(\Delta)\right] = \mathsf{E}\left[\exp\left\{-\hat{\theta}\Delta\right\}\right], \qquad \mathsf{Var}\left[\theta(\Delta)\right] = \mathsf{Var}\left[\exp\left\{-\hat{\theta}\Delta\right\}\right],$$

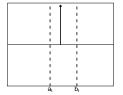
Georgi Dinolov (UCSC) March 11, 2019 60 / 66

Prior elicitation example IV

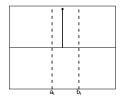
we obtain a system of two equations with two unknowns that can be solved numerically to find the values of $a_{\theta}(\Delta)$ and $b_{\theta}^2(\Delta)$ in terms of $\hat{a}_{\hat{\theta}}$, and Δ .

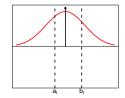
To elicit $\hat{a}_{\hat{\theta}}$ and $\hat{b}_{\hat{\theta}}^2$, recall that $\hat{\theta}$ is the inverse of the time scale of inertia for $\log(\hat{\sigma}_t)$ in the continuous-time formulation, which can be thought of as the characteristic time length, or unit of time, over which the process for the diffusion of $\log(\hat{\sigma}_t)$ "forgets" about an endogenous shock. The two hyper-parameters can be chosen so that the prior probability mass for $\hat{\theta}$ permits a reasonable range for the timescale of inertia.

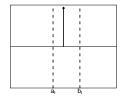


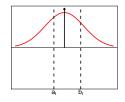


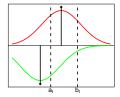


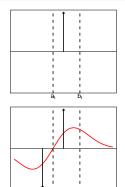


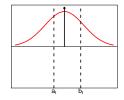


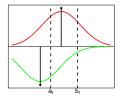


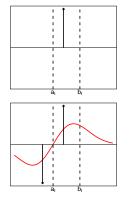


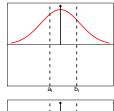


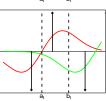


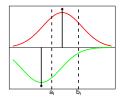


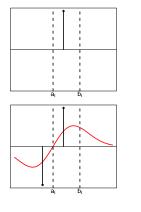


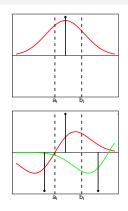


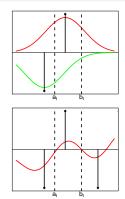


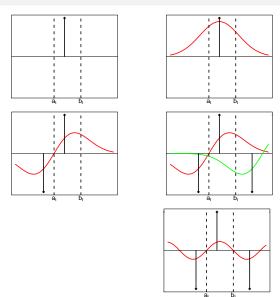


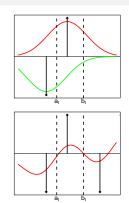








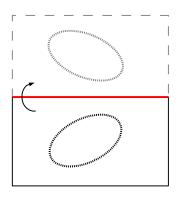


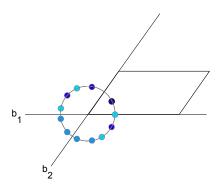




Solution by Method of Images: bivariate case

Method of Images does not work in generate.





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