

Nonlinear Filtering using Particles and Quadrature

Andreas Klöckner

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Outline

Nonlinear Filtering

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Nonlinear Filtering

Monte Carlo Filters

Particle Filters
Importance Sampling

Quadrature Filters

Error Estimates

Continuous-Time Filtering

- 1 Nonlinear Filtering
- 2 Monte Carlo Filters
 - Particle Filters
 - Importance Sampling
- 3 Quadrature Filters
 - Error Estimates
- 4 Continuous-Time Filtering

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The Nonlinear Filtering Problem

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Have:

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Have:

- A hidden *Markov process* \mathbf{x}_t for $t = 0, 1, 2, \dots$ with initial distribution $p(\mathbf{x}_0)$ and transition probability $p(\mathbf{x}_{t+1}|\mathbf{x}_t)$.

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- Conditionally independent *observations* \mathbf{y}_t for $t = 1, 2, 3, \dots$ characterized by $p(\mathbf{y}_t|\mathbf{x}_t)$.

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Want:

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- Conditionally independent *observations* \mathbf{y}_t for $t = 1, 2, 3, \dots$ characterized by $p(\mathbf{y}_t|\mathbf{x}_t)$.

Want:

- (An estimate of) The *posterior distribution* $p(\mathbf{x}_t|\mathbf{y}_{1:t})$.

Process and Observation, schematically

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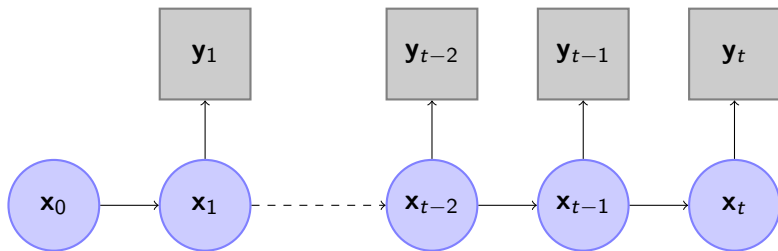
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Bayes' Formula

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Recall:

$$p(X|Y)$$

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Recall:

$$p(X|Y) = \frac{p(Y|X)p(X)}{p(Y)}$$

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Recall:

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So in our setting:

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So in our setting:

$$p(\mathbf{x}_{0:t}|\mathbf{y}_{1:t}) = \frac{p(\mathbf{y}_{1:t}|\mathbf{x}_{0:t})p(\mathbf{x}_{0:t})}{\int p(\mathbf{y}_{1:t}|\mathbf{x}_{0:t})p(\mathbf{x}_{0:t})d\mathbf{x}_{0:t}}$$

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Not usable for on-line processing.

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So in our setting:

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Not usable for on-line processing. Need a recursive algorithm.

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$$\frac{p(\mathbf{x}_{0:t} | \mathbf{y}_{1:t})}{p(\mathbf{x}_{0:t-1} | \mathbf{y}_{1:t-1})}$$

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$$= \frac{p(\mathbf{x}_{0:t} | \mathbf{y}_{1:t})}{p(\mathbf{x}_{0:t-1} | \mathbf{y}_{1:t-1})} \cdot \frac{p(\mathbf{y}_{1:t-1})}{p(\mathbf{y}_{1:t} | \mathbf{x}_{0:t}) p(\mathbf{x}_{0:t})} \cdot \frac{p(\mathbf{y}_{1:t-1})}{p(\mathbf{y}_{1:t-1} | \mathbf{x}_{0:t-1}) p(\mathbf{x}_{0:t-1})}$$

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$$\begin{aligned} & \frac{p(\mathbf{x}_{0:t}|\mathbf{y}_{1:t})}{p(\mathbf{x}_{0:t-1}|\mathbf{y}_{1:t-1})} \\ = & \frac{p(\mathbf{y}_{1:t}|\mathbf{x}_{0:t})p(\mathbf{x}_{0:t})}{p(\mathbf{y}_{1:t})} \cdot \frac{p(\mathbf{y}_{1:t-1})}{p(\mathbf{y}_{1:t-1}|\mathbf{x}_{0:t-1})p(\mathbf{x}_{0:t-1})} \\ \stackrel{\text{CI}}{=} & \frac{p(\mathbf{y}_t|\mathbf{x}_{0:t})p(\mathbf{x}_{0:t})}{p(\mathbf{y}_{1:t})} \cdot \frac{p(\mathbf{y}_{1:t-1})}{p(\mathbf{x}_{0:t-1})} \end{aligned}$$

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We can expand the denominator as

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We can expand the denominator as

$$p(\mathbf{y}_t|\mathbf{y}_{1:t-1}) = \int p(\mathbf{y}_t|\mathbf{x}_t)p(\mathbf{x}_t|\mathbf{y}_{1:t-1})d\mathbf{x}_t.$$

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Two drawbacks:

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Two drawbacks:

- Denominator not explicitly known

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We can expand the denominator as

$$p(\mathbf{y}_t|\mathbf{y}_{1:t-1}) = \int p(\mathbf{y}_t|\mathbf{x}_t)p(\mathbf{x}_t|\mathbf{y}_{1:t-1})d\mathbf{x}_t.$$

Two drawbacks:

- Denominator not explicitly known
- Don't care about the *joint* posterior—only want $p(\mathbf{x}_t|\mathbf{y}_{1:t})$

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$$p(\mathbf{x}_t | \mathbf{y}_{1:t})$$

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$$\stackrel{\text{JointUp}}{=} \int \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{x}_{t-1})}{p(\mathbf{y}_t | \mathbf{y}_{1:t-1})} p(\mathbf{x}_{0:t-1} | \mathbf{y}_{1:t-1}) d\mathbf{x}_{0:t-1}$$

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$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \int p(\mathbf{x}_{0:t} | \mathbf{y}_{1:t}) d\mathbf{x}_{0:t-1}$$

$$\stackrel{\text{JointUp}}{=} \int \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{x}_{t-1})}{p(\mathbf{y}_t | \mathbf{y}_{1:t-1})} p(\mathbf{x}_{0:t-1} | \mathbf{y}_{1:t-1}) d\mathbf{x}_{0:t-1}$$

$$= \frac{p(\mathbf{y}_t | \mathbf{x}_t)}{p(\mathbf{y}_t | \mathbf{y}_{1:t-1})} \int p(\mathbf{x}_t | \mathbf{x}_{t-1}) p(\mathbf{x}_{0:t-1} | \mathbf{y}_{1:t-1}) d\mathbf{x}_{0:t-1}$$

$$\stackrel{\text{Markov}}{=} \frac{p(\mathbf{y}_t | \mathbf{x}_t)}{p(\mathbf{y}_t | \mathbf{y}_{1:t-1})} \int p(\mathbf{x}_t | \mathbf{x}_{0:t-1}) p(\mathbf{x}_{0:t-1} | \mathbf{y}_{1:t-1}) d\mathbf{x}_{0:t-1}$$

$$= \frac{p(\mathbf{y}_t | \mathbf{x}_t)}{p(\mathbf{y}_t | \mathbf{y}_{1:t-1})} p(\mathbf{x}_t | \mathbf{y}_{1:t-1})$$

$$= \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})}{\int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t}$$

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$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})}{\int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t}$$

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Posterior $t - 1 \rightarrow$ Prediction t :

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$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})}{\int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t}$$

Posterior $t - 1 \rightarrow$ Prediction t :

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Prediction $t \rightarrow$ Posterior t :

$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})}{\int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t}$$

Posterior $t - 1 \rightarrow$ Prediction t :

Start with

$$p(\mathbf{x}_t) = \int p(\mathbf{x}_t | \mathbf{x}_{t-1}) p(\mathbf{x}_{t-1}) d\mathbf{x}_{t-1},$$

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Prediction $t \rightarrow$ Posterior t :

$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})}{\int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t}$$

Posterior $t - 1 \rightarrow$ Prediction t :

Start with

$$p(\mathbf{x}_t) = \int p(\mathbf{x}_t | \mathbf{x}_{t-1}) p(\mathbf{x}_{t-1}) d\mathbf{x}_{t-1},$$

and obtain

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Prediction $t \rightarrow$ Posterior t :

$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})}{\int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t}$$

Posterior $t - 1 \rightarrow$ Prediction t :

Start with

$$p(\mathbf{x}_t) = \int p(\mathbf{x}_t | \mathbf{x}_{t-1}) p(\mathbf{x}_{t-1}) d\mathbf{x}_{t-1},$$

and obtain

$$p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) = \int p(\mathbf{x}_t | \mathbf{x}_{t-1}) p(\mathbf{x}_{t-1} | \mathbf{y}_{1:t-1}) d\mathbf{x}_{t-1}.$$

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Our Example Process

Throughout this presentation, we will stick to the following model:

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Our Example Process

Throughout this presentation, we will stick to the following model:

$$x_t = \frac{1}{2}x_{t-1} + 25\frac{x_{t-1}}{1 + x_{t-1}^2}$$

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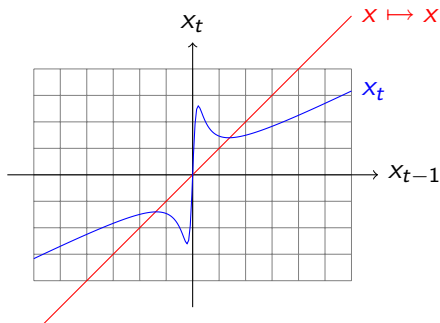
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Our Example Process

Throughout this presentation, we will stick to the following model:

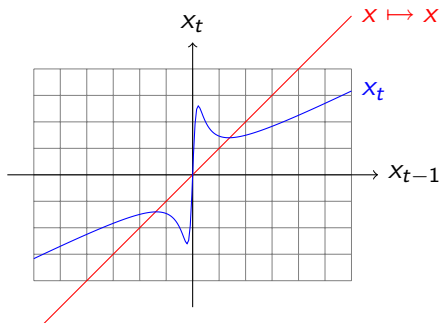
$$x_t = \frac{1}{2}x_{t-1} + 25 \frac{x_{t-1}}{1 + x_{t-1}^2}$$



Our Example Process

Throughout this presentation, we will stick to the following model:

$$x_t = \frac{1}{2}x_{t-1} + 25 \frac{x_{t-1}}{1 + x_{t-1}^2} + 8\cos(1.2t) + v_t,$$



Our Example Process

Throughout this presentation, we will stick to the following model:

$$x_t = \frac{1}{2}x_{t-1} + 25\frac{x_{t-1}}{1 + x_{t-1}^2} + 8\cos(1.2t) + v_t,$$
$$y_t = \frac{x_t^2}{20} + w_t,$$

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Throughout this presentation, we will stick to the following model:

$$x_t = \frac{1}{2}x_{t-1} + 25\frac{x_{t-1}}{1+x_{t-1}^2} + 8\cos(1.2t) + v_t,$$

$$y_t = \frac{x_t^2}{20} + w_t,$$

where $x_0 \sim \mathcal{N}(0, \sigma_1^2)$, $v_t \sim \mathcal{N}(0, \sigma_v^2)$, $w_t \sim \mathcal{N}(0, \sigma_w^2)$.

Particle Approximations

Assume a discrete approximation of the posterior density

$$\hat{p}(d\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{1}{N_t} \sum_{i=1}^{N_t} \delta_{\mathbf{x}_t^{(i)}}(d\mathbf{x}_t).$$

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Then insert that into the prediction formula:

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Then insert that into the prediction formula:

$$\hat{p}(d\mathbf{x}_{t+1} | \mathbf{y}_{1:t}) = \frac{1}{N_t} \sum_{i=1}^{N_t} p(d\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)})$$

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and the update formula:

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and the update formula:

$$\hat{p}(\mathbf{x}_{t+1} | \mathbf{y}_{1:t+1}) = \frac{1}{\tilde{C}_{t+1} \cdot N_t} \sum_{i=1}^{N_t} p(\mathbf{y}_{t+1} | \mathbf{x}_{t+1}) p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}).$$

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$$\hat{p}(d\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{1}{N_t} \sum_{i=1}^{N_t} \delta_{\mathbf{x}_t^{(i)}}(d\mathbf{x}_t).$$

Then insert that into the prediction formula:

$$\hat{p}(d\mathbf{x}_{t+1} | \mathbf{y}_{1:t}) = \frac{1}{N_t} \sum_{i=1}^{N_t} p(d\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)})$$

and the update formula:

$$\hat{p}(\mathbf{x}_{t+1} | \mathbf{y}_{1:t+1}) = \frac{1}{\tilde{c}_{t+1} \cdot N_t} \sum_{i=1}^{N_t} p(\mathbf{y}_{t+1} | \mathbf{x}_{t+1}) p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}).$$

Problem:

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Assume a discrete approximation of the posterior density

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and the update formula:

$$\hat{p}(\mathbf{x}_{t+1} | \mathbf{y}_{1:t+1}) = \frac{1}{\tilde{C}_{t+1} \cdot N_t} \sum_{i=1}^{N_t} p(\mathbf{y}_{t+1} | \mathbf{x}_{t+1}) p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}).$$

Problem: (Generally) Nontrivial to sample from!

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Importance Sampling lets you sample. . .

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Importance Sampling lets you sample...

- ... from an (almost) arbitrary density...

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Importance Sampling lets you sample...

- ... from an (almost) arbitrary density...
- ... that may only be known up to a proportionality constant.

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Importance Sampling lets you sample...

- ... from an (almost) arbitrary density...
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Recall our particle update formula:

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- ... that may only be known up to a proportionality constant.

Recall our particle update formula:

$$\hat{p}(\mathbf{x}_{t+1} | \mathbf{y}_{1:t+1}) = \frac{1}{\tilde{C}_{t+1} \cdot N_t} \sum_{i=1}^{N_t} p(\mathbf{y}_{t+1} | \mathbf{x}_{t+1}) p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}).$$

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That's exactly our situation.

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- . . . from an (almost) arbitrary density. . .
- . . . that may only be known up to a proportionality constant.

Recall our particle update formula:

$$\hat{p}(\mathbf{x}_{t+1} | \mathbf{y}_{1:t+1}) = \frac{1}{\tilde{C}_{t+1} \cdot N_t} \sum_{i=1}^{N_t} p(\mathbf{y}_{t+1} | \mathbf{x}_{t+1}) p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}).$$

That's exactly our situation.

Trick: Give each particle a *weight* in addition to its position.

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Suppose we have:

- $p(x) \propto q(x)$ not easy to sample from, q known exactly

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Suppose we have:

- $p(x) \propto q(x)$ not easy to sample from, q known exactly
- $\pi(x)$ a different density that is easy to sample from

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Suppose we have:

- $p(x) \propto q(x)$ not easy to sample from, q known exactly
- $\pi(x)$ a different density that is easy to sample from
–the *importance density*

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Suppose we have:

- $p(x) \propto q(x)$ not easy to sample from, q known exactly
- $\pi(x)$ a different density that is easy to sample from
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Then assign

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Suppose we have:

- $p(x) \propto q(x)$ not easy to sample from, q known exactly
- $\pi(x)$ a different density that is easy to sample from
–the *importance density*

Then assign

$$w^{(i)} \propto \frac{p(x^{(i)})}{\pi(x^{(i)})}$$

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Suppose we have:

- $p(x) \propto q(x)$ not easy to sample from, q known exactly
- $\pi(x)$ a different density that is easy to sample from
–the *importance density*

Then assign and normalize the weights

$$w^{(i)} \propto \frac{p(x^{(i)})}{\pi(x^{(i)})} \quad \tilde{w}^{(i)} := \frac{w^{(i)}}{\sum_j w^{(j)}}.$$

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Then assign and normalize the weights

$$w^{(i)} \propto \frac{p(x^{(i)})}{\pi(x^{(i)})} \quad \tilde{w}^{(i)} := \frac{w^{(i)}}{\sum_j w^{(j)}}.$$

Any unknown proportionality constant in p or w goes away in the normalization step.

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Now

$$p(x) \approx \hat{p}(x) := \sum_{i=1}^N \tilde{w}^{(i)} \delta(x - x^{(i)}),$$

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Now

$$p(x) \approx \hat{p}(x) := \sum_{i=1}^N \tilde{w}^{(i)} \delta(x - x^{(i)}),$$

assuming the particles $x^{(i)}$ have been drawn using $\pi(x)$.

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Why?

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$$p(x) \approx \hat{p}(x) := \sum_{i=1}^N \tilde{w}^{(i)} \delta(x - x^{(i)}),$$

assuming the particles $x^{(i)}$ have been drawn using $\pi(x)$.

Why?

A uniformly weighted distribution sampled from $p(x)$ is (approximately) the same as a distribution weighted with $p(x)/\pi(x)$ sampled from $\pi(x)$.

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Why?

A uniformly weighted distribution sampled from $p(x)$ is (approximately) the same as a distribution weighted with $p(x)/\pi(x)$ sampled from $\pi(x)$.

$$\int f(x) \frac{p(x)}{\pi(x)} \pi(x) dx = \int f(x) p(x) dx.$$

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Particle Framework + Importance Sampling = ?

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To maintain the recursive update property, demand

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Particle Framework + Importance Sampling = ?

To maintain the recursive update property, demand

$$\pi(\mathbf{x}_{0:t}|\mathbf{y}_{1:t}) = \pi(\mathbf{x}_{0:t-1}|\mathbf{y}_{1:t-1})\pi(\mathbf{x}_t|\mathbf{x}_{0:t-1}, \mathbf{y}_{1:t}).$$

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Recall the joint update equation

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Recall the joint update equation

$$p(\mathbf{x}_{0:t}|\mathbf{y}_{1:t}) \propto p(\mathbf{x}_{0:t-1}|\mathbf{y}_{1:t-1})p(\mathbf{y}_t|\mathbf{x}_t)p(\mathbf{x}_t|\mathbf{x}_{t-1}).$$

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This yields the *weight update equation*

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This yields the *weight update equation*

$$w_t^{(i)} \propto \frac{p(\mathbf{x}_{0:t}^{(i)}|\mathbf{y}_{1:t})}{\pi(\mathbf{x}_{0:t}^{(i)}|\mathbf{y}_{1:t})} =$$

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This yields the *weight update equation*

$$w_t^{(i)} \propto \frac{p(\mathbf{x}_{0:t}^{(i)}|\mathbf{y}_{1:t})}{\pi(\mathbf{x}_{0:t}^{(i)}|\mathbf{y}_{1:t})} = w_{t-1}^{(i)} \frac{p(\mathbf{y}_t|\mathbf{x}_t^{(i)})p(\mathbf{x}_t^{(i)}|\mathbf{x}_{t-1}^{(i)})}{\pi(\mathbf{x}_t^{(i)}|\mathbf{x}_{0:t-1}^{(i)}, \mathbf{y}_{1:t})}.$$

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Choosing a good $\pi(x)$ is extremely important.

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Choosing a good $\pi(x)$ is extremely important.

It can direct particles to interesting parts of the state space—or far away from there.

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Choosing a good $\pi(x)$ is extremely important.

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Goal: Keep the weights as uniform as possible.

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Choosing a good $\pi(x)$ is extremely important.

It can direct particles to interesting parts of the state space—or far away from there.

Goal: Keep the weights as uniform as possible.

Optimal choice:

$$\pi_{\text{opt}}(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t) = p(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t).$$

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Not usually easy to sample from → back to original problem.

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Popular choice:

$$\pi_{\text{prior}}(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t) := p(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)})$$

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Popular choice:

$$\pi_{\text{prior}}(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t) := p(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)})$$

Readily available and often Gaussian \rightarrow easy choice.

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Popular choice:

$$\pi_{\text{prior}}(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t) := p(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)})$$

Readily available and often Gaussian → easy choice.
Ignores observations → not always a good choice.

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Popular choice:

$$\pi_{\text{prior}}(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t) := p(\mathbf{x}_t | \mathbf{x}_{t-1}^{(i)})$$

Readily available and often Gaussian \rightarrow easy choice.

Ignores observations \rightarrow not always a good choice.

Weight update for π_{prior} :

$$w_t^{(i)} \propto w_{t-1}^{(i)} \frac{p(\mathbf{y}_t | \mathbf{x}_t^{(i)}) p(\mathbf{x}_t^{(i)} | \mathbf{x}_{t-1}^{(i)})}{\pi(\mathbf{x}_t^{(i)} | \mathbf{x}_{t-1}^{(i)}, \mathbf{y}_t)} = w_{t-1}^{(i)} p(\mathbf{y}_t | \mathbf{x}_t^{(i)}).$$

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\rightarrow Actually implementable algorithm.

SIS Implementation

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```
def sis(model, n=1000, max_time=20):
    particles = [model.sample_initial() for i in range(n)]
    particle_history = [particles]
    weights = [1/n for i in range(n)]
    weight_history = [weights]

    x = model.sample_initial()
    x_history = [x]

    for t in range(1, max_time):
        x = model.sample_transition(t, x)
        y = model.sample_observation(t, x)
        x_history.append(x)

        evolved_particles = [model.sample_transition(t, xp) for xp in particles]
        weights = [weight*model.observation_density(t, y, xtp)
                   for weight, xtp in zip(weights, evolved_particles)]
        weight_sum = sum(weights)
        weights = [weight/weight_sum for weight in weights]
        particle_history.append(particles)
        weight_history.append(weights)
```

SIS Results

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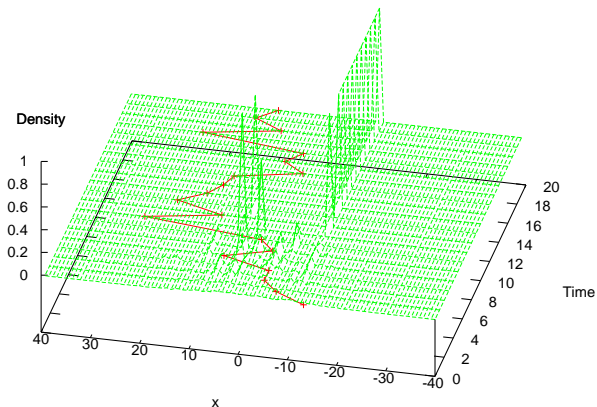
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SIS Method

trajectory —+—
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Sequential Importance Resampling

Pure SIS degenerates \rightarrow all weight on one particle.

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Pure SIS degenerates \rightarrow all weight on one particle.

Idea:

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Pure SIS degenerates→all weight on one particle.

Idea:

- Kill off particles with small weight

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Pure SIS degenerates→all weight on one particle.

Idea:

- Kill off particles with small weight
- Multiply the ones with high weight

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Pure SIS degenerates→all weight on one particle.

Idea:

- Kill off particles with small weight
- Multiply the ones with high weight
- Rely on process noise to scatter them

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How?

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Idea:

- Kill off particles with small weight
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How? Sample N times from

$$\hat{p}(x) := \sum_{i=1}^N \tilde{w}^{(i)} \delta(x - x^{(i)}).$$

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How? Sample N times from

$$\hat{p}(x) := \sum_{i=1}^N \tilde{w}^{(i)} \delta(x - x^{(i)}).$$

Particle descendant counts ξ satisfy

$$\xi \sim \text{Multinomial}(n, \tilde{w}^{(1)}, \dots, \tilde{w}^{(N)}).$$

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- Kill off particles with small weight
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Particle descendant counts ξ satisfy

$$\xi \sim \text{Multinomial}(n, \tilde{w}^{(1)}, \dots, \tilde{w}^{(N)}).$$

→ *Multinomial Resampling*

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```
def sir(model, n=1000, max_time=20):  
    particles = [model.sample_initial() for  
                i in range(n)]  
    particle_history = [particles]
```

```
x = model.sample_initial()  
x_history = [x]
```

```
for t in range(1, max_time):  
    x = model.sample_transition(t, x)  
    y = model.sample_observation(t, x)  
    x_history.append(x)
```

```
    evolved_particles = [model.  
                          sample_transition(t, xp) for xp  
                          in particles]  
    iweights = [model.  
                observation_density(t, y, xtp)  
                for xtp in evolved_particles]  
    particles = sample_discrete(  
        evolved_particles, iweights, n)  
    particle_history.append(particles)
```

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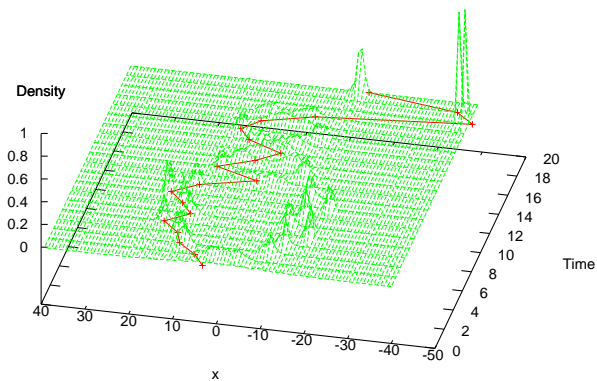
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SIR Method

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posterior - - - -



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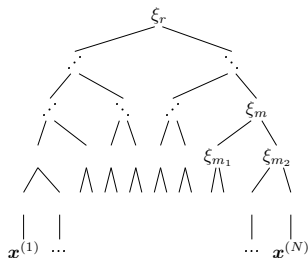
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Construct random variables $\xi_m \in \{0, \dots, N\}$ for each node m ,

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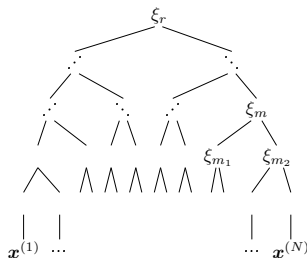
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Construct random variables $\xi_m \in \{0, \dots, N\}$ for each node m , with $\xi_m = \xi_{m_1} + \xi_{m_2}$

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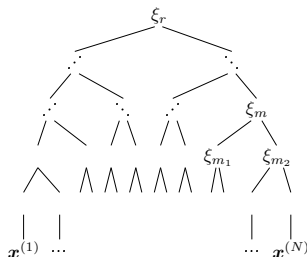
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Construct random variables $\xi_m \in \{0, \dots, N\}$ for each node m , with $\xi_m = \xi_{m_1} + \xi_{m_2}$ such that $E[\xi_m] = w_m$ and

$$\xi_m = \begin{cases} \lfloor w_m \rfloor & \text{with probability } 1 - w_m - \lfloor w_m \rfloor, \\ \lfloor w_m \rfloor + 1 & \text{with probability } w_m - \lfloor w_m \rfloor. \end{cases}$$

A Convergence Result

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Theorem (D. Crisan, 2001)

Let the transition kernel $p(\mathbf{x}_t|\mathbf{x}_{t-1})$ have the Feller property, that is for any bounded and continuous $f : \mathbb{R}^n \rightarrow \mathbb{R}$, the map

$$\mathbf{x}_{t-1} \mapsto \int f(x)p(d\mathbf{x}_t|\mathbf{x}_{t-1})$$

must be bounded and continuous. Then as the number of particles $N \rightarrow \infty$, the approximated posterior densities of the SIR method with multinomial resampling

$$\hat{p}^N(\mathbf{x}_t|\mathbf{y}_{0:t}) \rightarrow p(\mathbf{x}_t|\mathbf{y}_{0:t})$$

p -almost surely.

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Gauss-Legendre Quadrature

Idea: Cleverly reinterpret Gaussian Quadrature as a particle approximation.

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Idea: Cleverly reinterpret Gaussian Quadrature as a particle approximation.

Recall:

Gauss-Legendre Quadrature

Idea: Cleverly reinterpret Gaussian Quadrature as a particle approximation.

Recall:

$$\int_A^B f(x)dx \approx \sum_{i=1}^N f(x^{(i)})w^{(i)},$$

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Gauss-Legendre Quadrature

Idea: Cleverly reinterpret Gaussian Quadrature as a particle approximation.

Recall:

$$\int_A^B f(x)dx \approx \sum_{i=1}^N f(x^{(i)})w^{(i)},$$

with some quadrature weights $w^{(i)}$ and points $x^{(i)}$.

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Consider a probability density function p supported on (A, B) .

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Consider a probability density function p supported on (A, B) .

$$E[f(x)] = \int_A^B f(x)p(x)dx \approx \sum_{i=1}^N f(x^{(i)})p(x^{(i)})w^{(i)}.$$

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Consider a probability density function p supported on (A, B) .

$$E[f(x)] = \int_A^B f(x)p(x)dx \approx \sum_{i=1}^N f(x^{(i)})p(x^{(i)})w^{(i)}.$$

So view the quadrature points $x^{(i)}$ as the particles

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Idea: Cleverly reinterpret Gaussian Quadrature as a particle approximation.

Recall:

$$\int_A^B f(x)dx \approx \sum_{i=1}^N f(x^{(i)})w^{(i)},$$

with some quadrature weights $w^{(i)}$ and points $x^{(i)}$.

Consider a probability density function p supported on (A, B) .

$$E[f(x)] = \int_A^B f(x)p(x)dx \approx \sum_{i=1}^N f(x^{(i)})p(x^{(i)})w^{(i)}.$$

So view the quadrature points $x^{(i)}$ as the particles and

$$\tilde{p}(x^{(i)}) := p(x^{(i)})w^{(i)}$$

as the weights.

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Recall the Update Formula

$$p(\mathbf{x}_t | \mathbf{y}_{1:t}) = \frac{1}{C_t} \cdot p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1})$$

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$$C_t = \int p(\mathbf{y}_t | \mathbf{x}_t) p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) d\mathbf{x}_t.$$

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$$\tilde{p}(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t}) := \frac{1}{\hat{C}_t} \cdot p(\mathbf{y}_t | \mathbf{x}_t) w_t^{(i)} \underbrace{\hat{p}(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t-1})}_{\text{prediction}}.$$

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Compute this by finding

$$q^{(i)} := p(\mathbf{y}_t | \mathbf{x}_t) w_t^{(i)} \hat{p}(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t-1})$$

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$$q^{(i)} := p(\mathbf{y}_t | \mathbf{x}_t) w_t^{(i)} \hat{p}(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t-1})$$

$$\hat{C}_t = \sum_{i=1}^N q^{(i)},$$

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$$q^{(i)} := p(\mathbf{y}_t | \mathbf{x}_t) w_t^{(i)} \hat{p}(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t-1})$$

$$\hat{C}_t = \sum_{i=1}^N q^{(i)}, \quad \tilde{p}(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t}) = \frac{q^{(i)}}{\hat{C}_t}.$$

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$$\hat{p}(\mathbf{x}_{t+1}^{(i)}|\mathbf{y}_{1:t-1}) = \sum_{j=1}^N p(\mathbf{x}_{t+1}^{(i)}|\mathbf{x}_t^{(j)})\tilde{p}(\mathbf{x}_t^{(j)}|\mathbf{y}_{1:t})$$

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$\tilde{p}(\mathbf{x}_t^{(j)}|\mathbf{y}_{1:t})$ already contains quadrature weights.

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Note: $O(N^2)$ complexity.

Quadrature Filter Results

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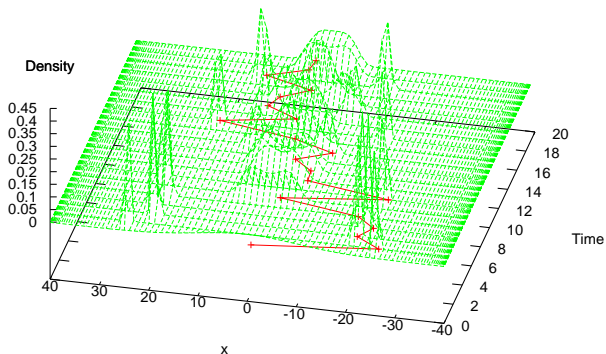
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trajectory —+—
posterior - - - - -



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- Deterministic, no sampling.

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- Deterministic, no sampling.
- Well-suited to comparing likelihoods

$$p(\mathbf{y}_{1:T}) = \prod_{t=1}^T p(\mathbf{y}_t | \mathbf{y}_{1:t-1}) = \prod_{t=1}^T C_t$$

of different models.

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of different models.

- Suffers heavily with high-dimensional statespaces: $O(N^2)$, $N = m^d$. (SIS: $O(N)$, SIR: $O(N \log N)$)

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- Requires far fewer particles than SMC for comparable accuracy.
- Explicit error estimates available.

Error Estimates: Setup

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Error in the log-likelihood of the observation:

$$\delta_t := \log\{\hat{p}(\mathbf{y}_{1:t})\} - \log\{p(\mathbf{y}_{1:t})\} = \sum_{s=1}^t \log \hat{C}_s - \sum_{s=1}^t \log C_s.$$

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Prediction error indicator:

$$\varepsilon(\mathbf{x}_{t+1}|\mathbf{y}_{1:t}) := \exp(\delta_t)\hat{p}(\mathbf{x}_{t+1}|\mathbf{y}_{1:t}) - p(\mathbf{x}_{t+1}|\mathbf{y}_{1:t}).$$

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Indicator ignores error in normalization \hat{C}_t

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Indicator ignores error in normalization $\hat{C}_t \rightarrow \exp(\delta_t)$ necessary.

Error Propagation Recursion

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A nice bit of algebra (see notes) shows

$$\begin{aligned} & \varepsilon(\mathbf{x}_{t+1}|\mathbf{y}_{1:t}) \\ = & C_t^{-1} \sum_i p(\mathbf{x}_{t+1}|\mathbf{x}_t^{(i)}) p(\mathbf{y}_t|\mathbf{x}_t^{(i)}) w_t^{(i)} \underline{\varepsilon(\mathbf{x}_t|\mathbf{y}_{1:t-1})} \\ & + \sum_i p(\mathbf{x}_{t+1}|\mathbf{x}_t^{(i)}) w_t^{(i)} p(\mathbf{x}_t^{(i)}|\mathbf{y}_{1:t}) - p(\mathbf{x}_{t+1}|\mathbf{y}_{1:t}). \end{aligned}$$

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Split error into

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Split error into

- Method error (First line)

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Split error into

- Method error (First line)
- Quadrature error (Second line) on *exact probabilities*

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- *Approximation Assumption*: Assume we have enough particles so that

Error Estimates: Assumptions

- *Approximation Assumption:* Assume we have enough particles so that

$$\left| \sum_j w^{(j)} p(\mathbf{x}_{t+1}^{(i)} | \mathbf{x}_t^{(j)}) p(\mathbf{x}_t^{(j)} | \mathbf{y}_{1:t}) - p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}) \right| \leq \varepsilon p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}).$$

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Note: Assumption is on quadrature of *exact* probabilities.

Error Estimates: Assumptions

- *Approximation Assumption*: Assume we have enough particles so that

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- *Induction Assumption*:

$$r_{t+1} \leq r_t(1 + \varepsilon) + \varepsilon,$$

Error Estimates: Assumptions

- *Approximation Assumption*: Assume we have enough particles so that

$$\left| \sum_j w^{(j)} p(\mathbf{x}_{t+1}^{(j)} | \mathbf{x}_t^{(j)}) p(\mathbf{x}_t^{(j)} | \mathbf{y}_{1:t}) - p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}) \right| \leq \varepsilon p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}).$$

Note: Assumption is on quadrature of *exact* probabilities.

- *Induction Assumption*:

$$r_{t+1} \leq r_t(1 + \varepsilon) + \varepsilon,$$

where r_t is the relative error at time t , i.e. the smallest positive real such that

$$|\varepsilon(\mathbf{x}_{t+1} | \mathbf{y}_{1:t})| \leq r_t p(\mathbf{x}_{t+1} | \mathbf{y}_{1:t}).$$

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$$\begin{aligned} & \varepsilon(\mathbf{x}_{t+1} | \mathbf{y}_{1:t}) \\ \stackrel{\text{IA}}{\leq} & |C_t^{-1} \sum_i p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}) p(\mathbf{y}_t | \mathbf{x}_t^{(i)}) w_t^{(i)} r_{t-1} p(\mathbf{x}_t | \mathbf{y}_{1:t-1})| \\ & + \varepsilon p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}) \end{aligned}$$

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$$\begin{aligned} & \varepsilon(\mathbf{x}_{t+1} | \mathbf{y}_{1:t}) \\ \stackrel{\text{IA}}{\leq} & \left| C_t^{-1} \sum_i p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}) p(\mathbf{y}_t | \mathbf{x}_t^{(i)}) w_t^{(i)} r_{t-1} p(\mathbf{x}_t | \mathbf{y}_{1:t-1}) \right| \\ & + \varepsilon p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}) \\ = & \left| r_{t-1} \sum_i p(\mathbf{x}_{t+1} | \mathbf{x}_t^{(i)}) w_t^{(i)} p(\mathbf{x}_t^{(i)} | \mathbf{y}_{1:t}) \right| \\ & + \varepsilon p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}) \\ \stackrel{\text{AA}}{\leq} & r_{t-1} |p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t})| (1 + \varepsilon) + \varepsilon p(\mathbf{x}_{t+1}^{(i)} | \mathbf{y}_{1:t}). \end{aligned}$$

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We showed

$$r_{t+1} \leq r_t(1 + \varepsilon) + \varepsilon.$$

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We showed

$$r_{t+1} \leq r_t(1 + \varepsilon) + \varepsilon.$$

This implies

$$r_t \leq (1 + \varepsilon)^t - 1.$$

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We showed

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→ Potentially *exponential* error growth.

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A similar proof shows

$$\delta_t \leq (t + 1) \log(1 + \varepsilon')$$

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A similar proof shows

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→ *Linear* error growth in the *log*-likelihood.

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Have:

- A *Markov diffusion process*

$$dx^i(t) = b^i(\mathbf{x}(t))dt + \sigma^{ij}(\mathbf{x}(t))dW^j(t)$$

where $\mathbf{x} = (x^1, \dots, x^d)$ and $\mathbf{x}(0) = \mathbf{x}_0$.

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- An *observation*

$$\mathbf{y}(t) = \int_0^t \mathbf{h}(\mathbf{x}(s))ds + \mathbf{V}(t).$$

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All coefficients are assumed smooth,

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All coefficients are assumed smooth, and \mathbf{x}_0 and the Wiener processes \mathbf{W} and \mathbf{V} are assumed to be independent.

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The optimal filter (= best mean-square estimate) for $f(\mathbf{x}(t))$ is

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The optimal filter (= best mean-square estimate) for $f(\mathbf{x}(t))$ is

$$\hat{f}(\mathbf{x}(t)) = \frac{\int_{\mathbb{R}^d} f(\mathbf{x}) u(t, \mathbf{x}) d\mathbf{x}}{\int_{\mathbb{R}^d} u(t, \mathbf{x}) d\mathbf{x}},$$

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where u is the *unnormalized filtering density*.

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where u is the *unnormalized filtering density*.

u obeys the *Zakai SPDE*

$$du(t, \mathbf{x}) = \mathcal{L}^* u(t, \mathbf{x}) dt + h(\mathbf{x}) u(t, \mathbf{x}) d\mathbf{y}(t),$$

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where

$$\mathcal{L}^* u := \frac{1}{2} \frac{\partial^2}{\partial x_i \partial x_j} ((\sigma \sigma^*)^{ij} u) - \frac{\partial}{\partial x_i} (b^i u).$$

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Sergey Lototsky's *spectral separating scheme* expresses u as an expansion into *Wick polynomials* ξ_α :

$$u(t, \mathbf{x}) = \sum_{\alpha} \frac{1}{\sqrt{\alpha!}} \varphi_{\alpha}(t, \mathbf{x}) \xi_{\alpha}(\mathbf{y}).$$

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$$u(t, \mathbf{x}) = \sum_{\alpha} \frac{1}{\sqrt{\alpha!}} \varphi_{\alpha}(t, \mathbf{x}) \xi_{\alpha}(\mathbf{y}).$$

- Coefficients φ_{α} satisfy a deterministic system of PDEs.
- The scheme separates
 - dependency on process parameters (φ_{α}) from
 - dependency on observation (ξ_{α}).
- φ_{α} can be precomputed.

Questions?

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