# Particle Filtering with Progressive Gaussian Approximations to the Optimal Importance Density

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#### The Plan

- A quick review of particle filtering and importance densities.
- The problem with the usual choices of importance density.
- The new progressive proposal method.
- How it relates to existing algorithms.
- Some simulation results.

# Particle Filtering

$$x_n \sim f(x_n|x_{n-1}) = \mathcal{N}(x_n|\phi(x_{n-1}), Q)$$
  
 $y_n \sim g(y_n|x_n) = \mathcal{N}(y_n|\psi(x_n), R)$   
 $x_1 \sim p(x_1) = \mathcal{N}(x_1|m_1, P_1),$ 

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• Particle filter approximates:

$$p(x_{1:n}|y_{1:n}) \propto f(x_n|x_{n-1})g(y_n|x_n)p(x_{1:n-1}|y_{1:n-1})$$

- Select an (n-1) particle and sample a new state,  $x_n^{(i)} \sim q(x_n|x_{n-1}^{(i)})$ .
- Update weight  $w_n^{(i)} = \frac{f(x_n^{(i)}|x_{n-1}^{(i)})g(y_n|x_n^{(i)})}{q(x_n^{(i)}|x_{n-1}^{(i)})}$ .



### Importance Densities

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Bootstrap filter

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Bootstrap filter

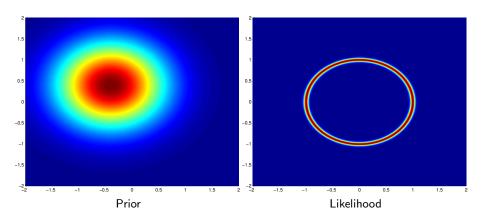
$$q(x_n|x_{n-1})=f(x_n|x_{n-1}).$$

Optimal Importance Density

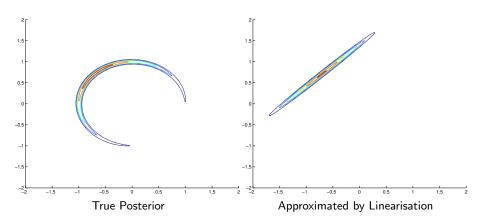
$$q(x_n|x_{n-1}) = \frac{f(x_n|x_{n-1})g(y_n|x_n)}{\int f(x_n|x_{n-1})g(y_n|x_n)dx_n}.$$

Approximate by linearisation or sigma-point approximation.

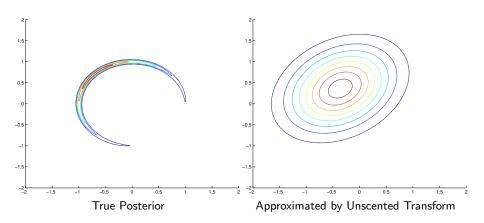
### An Example



### An Example - Linearisation



### An Example - Unscented Transform



### Progressive Principle

Introduce the observation gradually using a stretch of "pseudo-time",  $\lambda \in [0,1].$ 

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Smooth sequence of optimal importance densities:

$$\pi_{n,\lambda}(x_{n,\lambda}|x_{n-1}^{(j)}) \propto g(y_n|x_{n,\lambda})^{\lambda} f(x_{n,\lambda}|x_{n-1}^{(j)})$$

$$f(x_n|x_{n-1}) = \mathcal{N}(x_n | \phi(x_{n-1}), Q)$$
  
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Optimal importance density is analytically tractable.

$$\pi_{\lambda}(x_{\lambda}|x_{n-1}) = \mathcal{N}(x_{\lambda}|m_{\lambda}, P_{\lambda}),$$

$$P_{\lambda} = \left[ Q^{-1} + \lambda H^{T} R^{-1} H \right]^{-1}$$
  

$$m_{\lambda} = P_{\lambda} \left[ Q^{-1} \phi(x_{n-1}) + \lambda H^{T} R^{-1} y_{n} \right].$$

Any Gaussian random variable may be expressed as,

$$x = m + P^{\frac{1}{2}}z$$
$$z \sim \mathcal{N}(z|0,1).$$

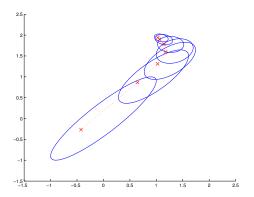
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Progress particle from  $\lambda_0$  to  $\lambda_1$  with a linear transformation.

$$x_{\lambda_1} = m_{\lambda_1} + P_{\lambda_1}^{\frac{1}{2}} P_{\lambda_0}^{-\frac{1}{2}} (x_{\lambda_0} - m_{\lambda_0}).$$

If 
$$x_{\lambda_0} \sim \pi_{\lambda_0}(x_{\lambda_0}|x_{n-1})$$
, then  $x_{\lambda_1} \sim \pi_{\lambda_1}(x_{\lambda_1}|x_{n-1})$ 



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Approximate the optimal importance density with a Gaussian at each point in pseudo-time.

$$\pi_{\lambda}(x_{\lambda}|x_{n-1}) \approx \mathcal{N}(x_{\lambda}|m_{\lambda}, P_{\lambda})$$

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Approximate the optimal importance density with a Gaussian at each point in pseudo-time.

$$\pi_{\lambda}(x_{\lambda}|x_{n-1}) \approx \mathcal{N}(x_{\lambda}|m_{\lambda}, P_{\lambda})$$

To update from  $\lambda_0$  to  $\lambda_1$ ,

$$\pi_{\lambda_{1}}(x|x_{n-1}) \propto \pi_{\lambda_{0}}(x|x_{n-1})g(y_{n}|x)^{\lambda_{1}-\lambda_{0}} \\ \propto \mathcal{N}(x|m_{\lambda_{0}}, P_{\lambda_{0}}) \mathcal{N}(y_{n}|\psi(x), R)^{\lambda_{1}-\lambda_{0}} \\ \propto \mathcal{N}(x|m_{\lambda_{0}}, P_{\lambda_{0}}) \mathcal{N}\left(y_{n} \middle| \psi(x), \frac{R}{\lambda_{1}-\lambda_{0}}\right).$$

Approximate with linearisation.

$$\hat{H}_{x_{\lambda_0}} = \left. \frac{\partial \psi}{\partial x} \right|_{x_{\lambda_0}}$$

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$$\mu_{\lambda_0} = \psi(x_{\lambda_0}) + \hat{H}_{x_{\lambda_0}}(m_{\lambda_0} - x_{\lambda_0})$$

$$\Sigma_{\lambda_0} = \hat{H}_{x_{\lambda_0}} P_{\lambda_0} \hat{H}_{x_{\lambda_0}}^T$$

$$C_{\lambda_0} = P_{\lambda_0} \hat{H}_{x_{\lambda_0}}^T$$

$$m_{\lambda_1} = m_{\lambda_0} + C_{\lambda_0} \left( \Sigma_{\lambda_0} + \frac{R}{\lambda_1 - \lambda_0} \right)^{-1} (y_n - \mu_{\lambda_0})$$

$$P_{\lambda_1} = P_{\lambda_0} - C_{\lambda_0} \left( \Sigma_{\lambda_0} + \frac{R}{\lambda_1 - \lambda_0} \right)^{-1} C_{\lambda_0}^T$$

### Weight Evolution

• Existing particle at pseudo time  $\lambda_0$ ,

$$\left\{x_{1:n-1}^{(i)}, x_{\lambda_0}^{(i)}\right\} \sim \eta_{\lambda_0}(x_{1:n-1}, x_{\lambda_0}).$$

• ... is replaced by new particle at  $\lambda_1$ ,

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• Standard change of variables formula,

$$\eta_{\lambda_1}(x_{1:n-1},x_{\lambda_1}) = \eta_{\lambda_0}(x_{1:n-1},x_{\lambda_0}) \times \left| \frac{\partial x_{\lambda_0}}{\partial x_{\lambda_1}} \right|.$$

### Weight Evolution

Hence weight update,

$$\begin{split} w_{\lambda_1} &= \frac{\tilde{\pi}_{\lambda_1}(x_{1:n-1}, x_{\lambda_1})}{\eta_{\lambda_1}(x_{1:n-1}, x_{\lambda_1})} \\ &= \frac{\tilde{\pi}_{\lambda_0}(x_{1:n-1}, x_{\lambda_0})}{\eta_{\lambda_0}(x_{1:n-1}, x_{\lambda_0})} \times \frac{\tilde{\pi}_{\lambda_1}(x_{1:n-1}, x_{\lambda_1})}{\tilde{\pi}_{\lambda_0}(x_{1:n-1}, x_{\lambda_0})} \times \left| \frac{\partial x_{\lambda_1}}{\partial x_{\lambda_0}} \right| \\ &\propto w_{\lambda_0} \times \frac{g(y_n|x_{\lambda_1})^{\lambda_1} f(x_{\lambda_1}|x_{n-1})}{g(y_n|x_{\lambda_0})^{\lambda_0} f(x_{\lambda_0}|x_{n-1})} \times \sqrt{\frac{|P_{\lambda_1}|}{|P_{\lambda_0}|}}. \end{split}$$

### Relationship to Other Methods

# Gradual introduction of likelihood underlies numerous existing particle methods,

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- N. Oudjane and C. Musso. Progressive correction for regularized particle filters.
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- F. Daum and J. Huang. Particle flow for nonlinear filters with log-homotopy.
   In Proceedings of SPIE, the International Society for Optical Engineering, pages 696918–1. Society of Photo-Optical Instrumentation Engineers, 2008
- Sebastian Reich. A dynamical systems framework for intermittent data assimilation. BIT Numerical Mathematics, 51:235–249, 2011

### Relationship to Other Methods

Distinguishing features of the new progressive proposal method:

- Particles moved deterministically.
- Move one particle at a time no intermediate interaction steps needed.
- Adaptive step-size control for each particle.

### Simulations - A Tracking Problem

$$x_n = \begin{bmatrix} p_n^T & v_n^T \end{bmatrix}^T$$

Near-constant velocity model.

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Near-constant velocity model.

$$y_n = \begin{bmatrix} \theta_n & r_n & h_n & s_n \end{bmatrix}^T$$
.

$$\theta_{n} = \arctan\left(\frac{p_{n,1}}{p_{n,2}}\right)$$

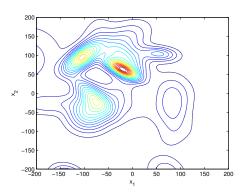
$$r_{n} = \sqrt{p_{n,1}^{2} + p_{n,3}^{2} + p_{n,3}^{2}}$$

$$h_{n} = p_{n,3} - T(p_{n,1}, p_{n,2})$$

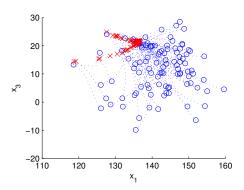
$$s_{n} = \frac{p_{n} \cdot v_{n}}{r_{n}},$$

# Simulations - Terrain Map





### Simulations - Results



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Algorithm	$N_F$	ESS	RMSE
Bootstrap Proposal	6000	1.0	78.6
Unscented Kalman Proposal	460	2.4	70.2
Gaussian Local Maximum Proposal	10	3.1	62.9
Progressive Proposal	180	56.4	22.3

### Summary

- The progressive proposal method samples from an effective approximation of the optimal importance density.
- Particles sampled from the transition density, then moved deterministically with a series of approximately optimal steps.
- Lower errors and better sample sizes when filtering with challenging nonlinear models.