

L9.1: An introduction to particle filtering

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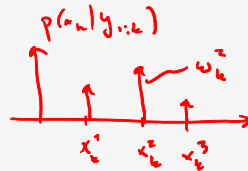
- Gaussian filtering is a useful technique to perform nonlinear filtering.
- **Limitations:** Gaussian filtering methods do not perform well when
 - the models are highly nonlinear,
 - when the posterior distribution is significantly non-Gaussian, e.g., a multimodal density.
- For such problems we need a different type of approximation to the posterior density!

Basic idea

- Use a non-parametric representation

$$\underline{p(\mathbf{x}_k | \mathbf{y}_{1:k})} \approx \sum_{i=1}^N \underline{w_k^{(i)}} \delta(\mathbf{x}_k - \underline{\mathbf{x}_k^{(i)}})$$

where $\mathbf{x}_k^{(i)}$ are particles and $w_k^{(i)}$ are associated weights.



- Filtering is (essentially) performed by
 1. propagating $\underline{\mathbf{x}_{k-1}^{(i)}} \rightarrow \underline{\mathbf{x}_k^{(i)}}$ over time,
 2. updating the weights, $\underline{w_k^{(i)}}$.

- Basic version: $\underline{\mathbf{x}_k^{(i)}} \sim \underline{p(\mathbf{x}_k | \mathbf{x}_{k-1}^{(i)})}$, $\underline{w_k^{(i)}} \propto \underline{w_{k-1}^{(i)}} \underline{p(\mathbf{y}_k | \mathbf{x}_k^{(i)})}$.

After this lecture you should be able to

- explain the concepts of Monte Carlo sampling and importance sampling,
- describe what particle degeneracy is and why resampling is useful,
- and implement a particle filter.

L9.2: Monte Carlo (MC) approximations and Importance Sampling (IS)

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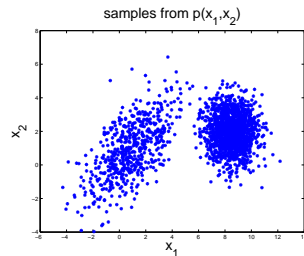
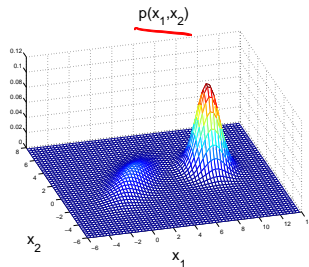
Monte Carlo approximations

Two perspectives on Monte Carlo approximation

Given independent samples $\mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(N)} \sim p(\mathbf{x})$ we can approximate

$$\mathbb{E}[\mathbf{g}(\mathbf{x})] \approx \frac{1}{N} \sum_{i=1}^N \mathbf{g}(\mathbf{x}^{(i)}) \quad (1)$$

$$p(\mathbf{x}) \approx \frac{1}{N} \sum_{i=1}^N \delta(\mathbf{x} - \mathbf{x}^{(i)}) \quad (2)$$



Remarks on Monte Carlo approximations:

- non-parametric approximation to $p(\mathbf{x})$.
- approximate all kinds of densities, $p(\mathbf{x})$. *Very flexible!*
- does not suffer from the curse of dimensionality, e.g., $\hat{\mu}$

$$\text{Cov}(\hat{\mu}) = \text{Cov} \left(\underbrace{\frac{1}{N} \sum_{i=1}^N \mathbf{x}^{(i)}}_{\hat{\mu}} \right) = \frac{1}{N} \text{Cov}(\underline{\mathbf{x}})$$

independently on $\dim(\mathbf{x})$!

- *Weakness:* it is often difficult to generate samples from $p(\mathbf{x})$.

Importance sampling

What can we do when it is difficult to sample from $p(\mathbf{x})$?

Importance sampling

- Generate samples, $\mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(N)}$, from a proposal density $q(\mathbf{x})$:

$$\mathbb{E}_{\underline{p(\mathbf{x})}}[\underline{\mathbf{g}(\mathbf{x})}] = \int \underbrace{\mathbf{g}(\mathbf{x}) \cdot \frac{p(\mathbf{x})}{q(\mathbf{x})}}_{\tilde{\mathbf{g}}(\mathbf{x})} d\mathbf{x} \approx \frac{1}{N} \sum_{i=1}^N \mathbf{g}(\mathbf{x}^{(i)}) \frac{p(\mathbf{x}^{(i)})}{q(\mathbf{x}^{(i)})}$$

$$\propto \sum_{i=1}^N \underbrace{\frac{p(\mathbf{x}^{(i)})}{q(\mathbf{x}^{(i)})}}_{\tilde{w}(i)} \cdot \mathbf{g}(\mathbf{x}^{(i)})$$

Importance sampling approximation to $p(\mathbf{x})$

- Generate samples, $\mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(N)}$, from $q(\mathbf{x})$ and set

$$\underline{p(\mathbf{x})} \approx \sum_{i=1}^N \underline{w^{(i)}} \underline{\delta(\mathbf{x} - \mathbf{x}^{(i)})}$$

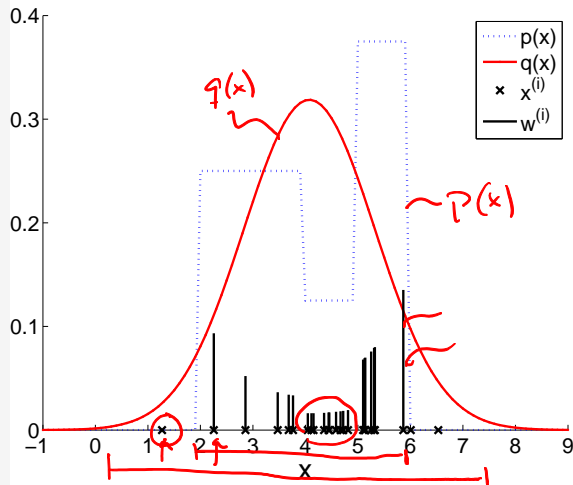
where

$$\underline{w^{(i)}} = \frac{\tilde{w}^{(i)}}{\sum_{n=1}^N \tilde{w}^{(n)}} \quad \text{and} \quad \tilde{w}^{(i)} = \frac{p(\mathbf{x}^{(i)})}{q(\mathbf{x}^{(i)})}.$$

- Importance sampling is a flexible and powerful tool.
- It can perform very well as long as:
 1. it is easy to sample from $q(\mathbf{x})$,
 2. the support of $q(\mathbf{x})$ contains the support of $p(\mathbf{x})$,
 3. $q(\mathbf{x})$ is “similar” to $p(\mathbf{x})$.

Example – Importance sampling

- Approximate $p(x)$ using N independent samples from $q(x) = \mathcal{N}(x; 4, 1.5^2)$.



L9.3: Sequential Importance Sampling (SIS)

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- **Objective:** to recursively and accurately approximate the filtering density, $p(\mathbf{x}_k | \mathbf{y}_{1:k})$.

- **Assumption:** both the motion and measurement models

$$p(\mathbf{x}_k | \mathbf{x}_{k-1}) \quad \text{and} \quad p(\mathbf{y}_k | \mathbf{x}_k)$$

can be easily evaluated point-wise.

- A common example is

$$\mathbf{x}_k = \underline{f(\mathbf{x}_{k-1})} + \underline{\mathbf{q}_{k-1}}, \quad \mathbf{q}_{k-1} \sim \mathcal{N}(\mathbf{0}, \mathbf{Q}_{k-1})$$

$$\mathbf{y}_k = \underline{h(\mathbf{x}_k)} + \underline{\mathbf{r}_k}, \quad \mathbf{r}_k \sim \mathcal{N}(\mathbf{0}, \mathbf{R}_k),$$

where, e.g., $p(\mathbf{x}_k | \mathbf{x}_{k-1}) = \mathcal{N}(\mathbf{x}_k; f(\mathbf{x}_{k-1}), \mathbf{Q}_{k-1})$ is generally easy to evaluate for any values of \mathbf{x}_k and \mathbf{x}_{k-1} .

- **Particle filters** are also known as sequential importance resampling or *sequential Monte Carlo*.
- The basis of these methods is an algorithm called sequential importance sampling (SIS).

Standard SIS algorithm

- For $i = 1, \dots, N$ and at each time k :
 - Draw $\mathbf{x}_k^{(i)} \sim q(\mathbf{x}_k | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)$.
 - Compute weights

$$\underline{w_k^{(i)}} \propto \underline{w_{k-1}^{(i)}} \frac{\underline{p(\mathbf{y}_k | \mathbf{x}_k^{(i)}) p(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)})}}{\underline{q(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)}}.$$

- Normalize the weights.

- We then approximate $p(\mathbf{x}_k | \mathbf{y}_{1:k}) \approx \sum_{i=1}^N w_k^{(i)} \delta(\mathbf{x}_k - \mathbf{x}_k^{(i)})$.

- Assuming that we describe our posterior using the following approximation $p(\mathbf{x}_k | \mathbf{y}_{1:k}) \approx \sum_{i=1}^N w_k^{(i)} \delta(\mathbf{x}_k - \mathbf{x}_k^{(i)})$. What is then the MMSE estimate of \mathbf{x}_k ?

- $\hat{\mathbf{x}}_k = \sum_i^N w_k^{(i)} \mathbf{x}_k^{(i)}$

- $\hat{\mathbf{x}}_k = \mathbf{x}_k^{(j)}$, where $j = \arg \max_i w_k^{(i)}$

- $\hat{\mathbf{x}}_k = \frac{1}{N} \sum_i^N \mathbf{x}_k^{(i)}$

$$\begin{aligned} \hat{\mathbf{x}}_k &= E\{\mathbf{x}_k | \mathbf{y}_k\} = \int \mathbf{x}_k \sum_{i=1}^N w_k^{(i)} \delta(\mathbf{x}_k - \mathbf{x}_k^{(i)}) d\mathbf{x}_k \\ &= \sum_{i=1}^N w_k^{(i)} \mathbf{x}_k^{(i)} \end{aligned}$$

- It is not possible to calculate a MMSE estimate from this approximation.

Derivation - Basic strategy

Recursively at time $k = 1, 2, \dots$

1. Draw particles

$$\underline{\mathbf{x}_{0:k}^{(i)}} \sim \underline{q(\mathbf{x}_{0:k} | \mathbf{y}_{1:k})}$$

2. Update weights

$$\underline{w_k^{(i)} \propto \frac{p(\mathbf{x}_{0:k}^{(i)} | \mathbf{y}_{1:k})}{q(\mathbf{x}_{0:k}^{(i)} | \mathbf{y}_{1:k})}}$$

Comments on drawing particles:

- Let us **assume that**

$$\underline{q(\mathbf{x}_{0:k} | \mathbf{y}_{1:k})} = \underline{q(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k)} \underline{q(\mathbf{x}_{0:k-1} | \mathbf{y}_{1:k-1})}.$$

- we generate $\underline{\mathbf{x}_{0:k-1}^{(i)} \sim q(\mathbf{x}_{0:k-1} | \mathbf{y}_{1:k-1})}$ at time $k-1$,
- it is sufficient to generate $\underline{\mathbf{x}_k^{(i)} \sim q(\mathbf{x}_k | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)}$ and append that to $\underline{\mathbf{x}_{1:k-1}^{(i)}}$.

- It remains to derive the expression for the weights:

$$w_k^{(i)} \propto \frac{p(\mathbf{x}_{0:k}^{(i)} | \mathbf{y}_{1:k})}{\underline{q(\mathbf{x}_{0:k}^{(i)} | \mathbf{y}_{1:k})}} \propto p(\mathbf{x}_{0:k-1}^{(i)}, \mathbf{x}_k^{(i)} | \mathbf{y}_{1:k-1}) = p(\mathbf{y}_k | \mathbf{x}_k^{(i)}) p(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)}) p(\mathbf{x}_{0:k-1}^{(i)} | \mathbf{y}_{1:k-1})$$

$$\propto \frac{p(\mathbf{y}_k | \mathbf{x}_k^{(i)}) p(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)})}{q(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)} \underbrace{\frac{p(\mathbf{x}_{0:k-1}^{(i)} | \mathbf{y}_{1:k-1})}{q(\mathbf{x}_{0:k-1}^{(i)} | \mathbf{y}_{1:k-1})}}_{w_{k-1}^{(i)}}$$

$$\propto w_{k-1}^{(i)} \frac{p(\mathbf{y}_k | \mathbf{x}_k^{(i)}) p(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)})}{q(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)}$$

- We have thus derived the SIS algorithm:

Standard SIS algorithm

- For $i = 1, \dots, N$ and at each time k :

- Draw $\mathbf{x}_k^{(i)} \sim q(\mathbf{x}_k | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)$.

- Compute weights

$$\underline{w_k^{(i)} \propto w_{k-1}^{(i)} \frac{p(\mathbf{y}_k | \mathbf{x}_k^{(i)}) p(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)})}{q(\mathbf{x}_k^{(i)} | \mathbf{x}_{k-1}^{(i)}, \mathbf{y}_k)}}.$$

- Normalize the weights.

- A simple choice of **importance density** is

$$q(\mathbf{x}_k | \mathbf{x}_{k-1}, \cancel{\mathbf{y}_k}) = \underline{p(\mathbf{x}_k | \mathbf{x}_{k-1})}$$

for which $w_k^{(i)} \propto w_{k-1}^{(i)} p(\mathbf{y}_k | \mathbf{x}_k^{(i)})$.

Example - Nonlinear filter benchmark

- The following is a common benchmark for nonlinear filters

$$x_k = \frac{x_{k-1}}{2} + \frac{25x_{k-1}}{1 + x_{k-1}^2} + 8 \cos(1.2k) + q_{k-1}$$

$$y_k = \frac{x_k^2}{20} + r_k$$

where $q_{k-1} \sim \mathcal{N}(0, 10)$ and $r_k \sim \mathcal{N}(0, 1)$.

- Let us see how the above filter performs on this challenging problem!

L9.4: Sequential Importance Resampling (SIR)

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- One can show that all SIS filters suffer from **degeneracy**:
after a few time steps all but one particle will have negligible weight.
- Consequences of degeneracy:
 - the filter believes that it knows \mathbf{x}_k exactly,
 - we obtain very poor state estimates,
 - most of our calculations are wasted on insignificant particles.

These are very serious drawbacks!

- A key technique to improve performance is **resampling**.

- **Challenge:** we have $p(\mathbf{x}_k | \mathbf{y}_{1:k}) \approx \sum_{i=1}^N w_k^{(i)} \delta(\mathbf{x}_k - \mathbf{x}_k^{(i)})$ where most weights $w_k^{(i)}$ are very small.

Idea: use Monte Carlo sampling

- Generate independent samples $\tilde{\mathbf{x}}_k^{(1)}, \dots, \tilde{\mathbf{x}}_k^{(N)}$ from $p(\mathbf{x}_k | \mathbf{y}_{1:k})$ and set

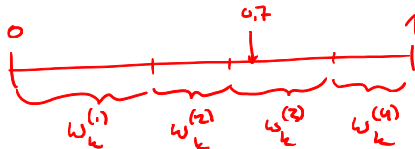
$$p(\mathbf{x}_k | \mathbf{y}_{1:k}) \approx \sum_{i=1}^N \frac{1}{N} \delta(\mathbf{x}_k - \tilde{\mathbf{x}}_k^{(i)}).$$

- After resampling we get
 - equal weights (they are all $1/N$),
 - multiple copies of high probability particles.

Resampling algorithm

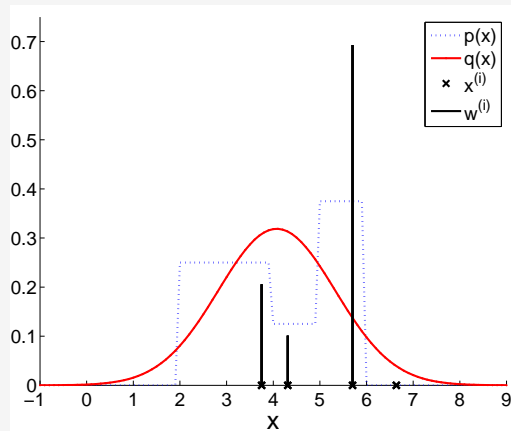
- 1) Draw N samples with replacement from $\mathbf{x}_k^{(1)}, \mathbf{x}_k^{(2)}, \dots, \mathbf{x}_k^{(N)}$, where the probability of selecting $\mathbf{x}_k^{(i)}$ is $w_k^{(i)}$.
- 2) Replace the old sample set with the new one and set all weights to $1/N$.

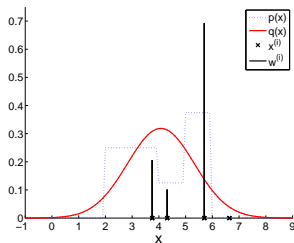
- A few remarks:
 - We use $\mathbf{x}_k^{(i)}$ and $w_k^{(i)}$ to denote the particles and their weights also after resampling.
 - We can use samples from the uniform distribution, $\text{unif}[0, 1]$, to draw samples from the discrete distribution $p(\mathbf{x}_k | \mathbf{y}_{1:k})$.



Self-assessment – Resampling

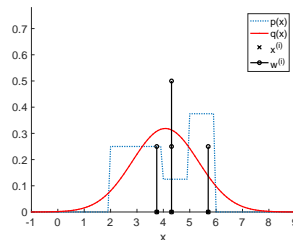
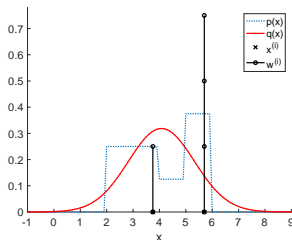
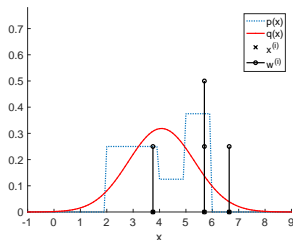
- Perform resampling on the density to the right and illustrate the result.
- Assume that the numbers 0.65, 0.03, 0.84 and 0.93 are drawn from $\text{unif}[0, 1]$.



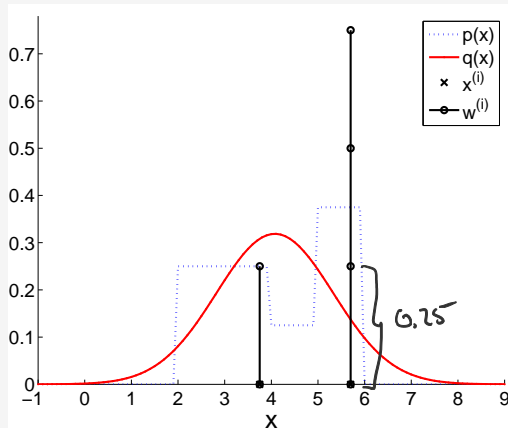


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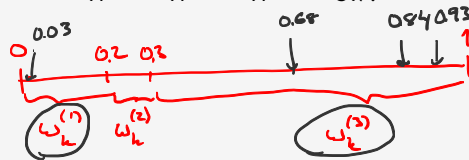
• Choose the figure below that illustrates the resampled particles:



Self-assessment – Solution



- If particles were ordered in ascending order, $x^{(1)} < \dots < x^{(4)}$, resampling gives $x^{(1)} = 3.8$ and $x^{(2)} = x^{(3)} = x^{(4)} = 5.7$.



- Resampling costs some calculations and introduces some errors, but improves performance immensely over time.
- An estimate for the *effective number of particles* is

$$N_{eff} = \frac{1}{\sum_{i=1}^N \left(w_k^{(i)}\right)^2}.$$

- Many algorithms only resample when N_{eff} is below some threshold, e.g., $N/4$.

L9.5: Choice of importance distribution

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- A carefully selected importance distribution, $q(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k)$, can slow down the degeneracy and improve performance.

Intuition: if most particles are placed in "high probability regions" there is less need to get rid of useless particles.

Optimal importance density

- The optimal importance density is

$$q(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k) = p(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k).$$

- Unfortunately, in most nonlinear settings, $p(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k)$, is difficult to both draw samples from and to evaluate.

- We can approximate $p(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k)$ using, e.g., linearization.
- The most common choice is still, $q(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{y}_k) = p(\mathbf{x}_k | \mathbf{x}_{k-1})$, and the bootstrap algorithm.
- **Note:** if we resample at every time step, we get $w_k^{(i)} \propto p(\mathbf{y}_k | \mathbf{x}_k^{(i)})$ since $w_{k-1}^{(i)} = 1/N \ \forall i$ after resampling.
- **Note 2:** the Auxiliary PF (APF) is variation of the SIR algorithm that makes use of \mathbf{y}_k .

The bootstrap PF

At each time k :

- Draw $\mathbf{x}_k^{(i)} \sim p(\mathbf{x}_k | \mathbf{x}_{k-1}^{(i)})$, for $i = 1, \dots, N$.
- Calculate $w_k^{(i)} \propto w_{k-1}^{(i)} p(\mathbf{y}_k | \mathbf{x}_k^{(i)})$ and normalize to 1.
- Resample.

- Particle filters (PFs) can handle highly nonlinear and non-Gaussian systems.
- Particle filters are asymptotically exact as you increase N .
- The complexity is roughly $O(N)$ but the gain in performance flattens out as you increase N .
- Unfortunately, PFs suffer from the curse of dimensionality and are intractable in higher dimensions.

- The output from a PF is an approximation

$$\underline{p(\mathbf{x}_k | \mathbf{y}_{1:k}) \approx \sum_{i=1}^N w_k^{(i)} \delta(\mathbf{x}_k - \mathbf{x}_k^{(i)})}$$

$$\Rightarrow Pr\{\mathbf{x}_k = \mathbf{x}' | \mathbf{y}_{1:k}\} = \begin{cases} \omega_k^{(i)} & \text{if } \mathbf{x}' = \mathbf{x}_k^{(i)} \\ 0 & \text{otherwise} \end{cases}$$

which implies that

$$\mathbb{E} [\mathbf{g}(\mathbf{x}_k) | \mathbf{y}_{1:k}] \approx \sum_{i=1}^N w_k^{(i)} \mathbf{g}(\mathbf{x}_k^{(i)}).$$

L9.6: Rao-Blackwellized Particle Filter

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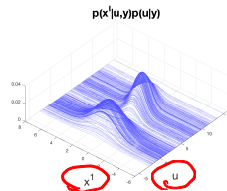
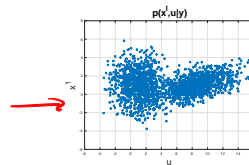
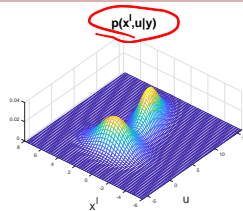
- Background:
 - particle filters are intractable in high dimensions.
 - many systems are linear in some dimensions.

• Idea 1: “combine a particle filter for the nonlinear states with a Kalman filter for the linear states”.

• Idea 2: If $\mathbf{x}_k = \begin{bmatrix} \mathbf{x}_k^l \\ \mathbf{u}_k \end{bmatrix}$ where \mathbf{x}_k^l and \mathbf{u}_k are the linear and nonlinear states:

$$p(\mathbf{x}_k^l, \mathbf{u}_{1:k} | \mathbf{y}_{1:k}) = p(\mathbf{x}_k^l | \underbrace{\mathbf{u}_{0:k}}_{\text{Gaussian}}, \underbrace{\mathbf{y}_{1:k}}_{\text{Particle filter}}) p(\mathbf{u}_{0:k} | \mathbf{y}_{1:k})$$

Gaussian Particle filter



- Assuming we have $\mathbf{x}_k = \begin{bmatrix} \mathbf{x}_k^l \\ \mathbf{u}_k \end{bmatrix}$, Rao-Blackwellized particle filters are often used for models on the form

$$\mathbf{x}_k^l = \underline{f_{k-1}^l(\mathbf{u}_{k-1})} + \underline{\mathbf{A}_{k-1}^l(\mathbf{u}_{k-1})} \underline{\mathbf{x}_{k-1}^l} + \mathbf{q}_{k-1}^l$$

$$\mathbf{u}_k = \underline{f_{k-1}^u(\mathbf{u}_{k-1})} + \underline{\mathbf{A}_{k-1}^u(\mathbf{u}_{k-1})} \underline{\mathbf{x}_{k-1}^l} + \mathbf{q}_{k-1}^u$$

$$\mathbf{y}_k = \underline{h_k(\mathbf{u}_k)} + \underline{\mathbf{H}_k(\mathbf{u}_k)} \underline{\mathbf{x}_k^l} + \mathbf{r}_k$$

where all the noises are Gaussian.

Bearing only tracking

- Bearing only tracking with a constant velocity motion in 2D.

What is \mathbf{x}_k^l , \mathbf{u}_k^l and \mathbf{y}_k in this example?

- \mathbf{x}_k^l : position, \mathbf{u}_k^l : velocity, \mathbf{y}_k : bearing to target
- \mathbf{x}_k^l : velocity, \mathbf{u}_k^l : position, \mathbf{y}_k : bearing to target
- \mathbf{x}_k^l : velocity, \mathbf{u}_k^l : bearing to target, \mathbf{y}_k : position
- \mathbf{x}_k^l : position, \mathbf{u}_k^l : bearing to target, \mathbf{y}_k : bearing to target

Bearing only tracking – system models

- Let us denote our state vector $\mathbf{x}_k = [\underbrace{x_k^1, x_k^2}_{u_k}, \underbrace{\dot{x}_k^1, \dot{x}_k^2}_{x_k^l}]^T$, the system models can then be written as:

$$\begin{aligned} x_k^l &= \begin{bmatrix} \dot{x}_k^1 \\ \dot{x}_k^2 \end{bmatrix} = \begin{bmatrix} \dot{x}_{k-1}^1 \\ \dot{x}_{k-1}^2 \end{bmatrix} + \mathbf{q}_{k-1}^l = x_{k-1}^l - q_{k-1}^l \\ u_k &= \begin{bmatrix} x_k^1 \\ x_k^2 \end{bmatrix} = \begin{bmatrix} x_{k-1}^1 \\ x_{k-1}^2 \end{bmatrix} + T \begin{bmatrix} \dot{x}_{k-1}^1 \\ \dot{x}_{k-1}^2 \end{bmatrix} + \mathbf{q}_{k-1}^u = u_{k-1} + T \cdot x_{k-1}^l + q_{k-1}^u \end{aligned}$$

$$\mathbf{y}_k = \text{atan}_2(x_k^2, x_k^1) + \mathbf{r}_k$$

$$\text{where } \mathbf{r}_k \sim \mathcal{N}(0, \sigma_r^2) \text{ and } \mathbf{q}_k = \begin{bmatrix} \mathbf{q}_k^u \\ \mathbf{q}_k^l \end{bmatrix} \sim \mathcal{N}\left(\mathbf{0}, \begin{bmatrix} \frac{T^2}{2} \mathbf{I} \\ \Pi \end{bmatrix} \begin{bmatrix} \sigma_q^2 & 0 \\ 0 & \sigma_q^2 \end{bmatrix} \begin{bmatrix} \frac{T^2}{2} \mathbf{I} \\ \Pi \end{bmatrix}^T\right)$$

- One recursion of the Rao-Blackwellized particle filter contains five steps:

$p(\mathbf{x}_{k-1}^l \mathbf{u}_{0:k-1}, \mathbf{y}_{1:k-1}) p(\mathbf{u}_{0:k-1} \mathbf{y}_{1:k-1}) \rightarrow p(\mathbf{x}_k^l \mathbf{u}_{0:k}, \mathbf{y}_{1:k}) p(\mathbf{u}_{0:k} \mathbf{y}_{1:k})$	
1) PF-pred:	$p(\mathbf{u}_{1:k-1} \mathbf{y}_{1:k-1}) \rightarrow p(\mathbf{u}_{1:k} \mathbf{y}_{1:k-1})$
2) KF, dyn. upd.:	$p(\mathbf{x}_{k-1}^l \mathbf{u}_{1:k-1}, \mathbf{y}_{1:k-1}) \rightarrow p(\mathbf{x}_{k-1}^l \mathbf{u}_{1:k}, \mathbf{y}_{1:k-1})$
3) KF-pred:	$p(\mathbf{x}_{k-1}^l \mathbf{u}_{1:k-1}, \mathbf{y}_{1:k-1}) \rightarrow p(\mathbf{x}_k^l \mathbf{u}_{1:k-1}, \mathbf{y}_{1:k-1})$
4) PF, update:	$p(\mathbf{u}_{1:k} \mathbf{y}_{1:k-1}) \rightarrow p(\mathbf{u}_{1:k} \mathbf{y}_{1:k})$
5) KF, meas. upd.:	$p(\mathbf{x}_k^l \mathbf{u}_{1:k}, \mathbf{y}_{1:k-1}) \rightarrow p(\mathbf{x}_k^l \mathbf{u}_{1:k}, \mathbf{y}_{1:k})$

$$\mathbf{u}_k^{(i)} = \mathbf{u}_{k-1}^{(i)} + \mathbf{T} \cdot \mathbf{x}_k^l + \mathbf{q}_{k-1}^u$$

$$\mathbf{u}_k^{(i)} - \mathbf{u}_{k-1}^{(i)} = \mathbf{T} \cdot \mathbf{x}_k^l + \mathbf{q}_{k-1}^u$$

$$\mathbf{x}_k^l = \mathbf{x}_{k-1}^l + \mathbf{q}_{k-1}^l$$

$$\omega_k^{(i)} \propto \omega_{k-1}^{(i)} \cdot \mathcal{P}(y_k | u_k)$$

- Note:**

- Step 2) makes use of the motion model for \mathbf{u}_k to update \mathbf{x}_{k-1}^l .
- The linear states are marginalized from step 1) and 4), similarly to how we normally handle noise.

Bearing only tracking – system models

- Let us denote our state vector $\mathbf{x}_k = [x_k^1, x_k^2, \dot{x}_k^1, \dot{x}_k^2]^T$, the system models can then be written as:

$$\begin{bmatrix} \dot{x}_k^1 \\ \dot{x}_k^2 \end{bmatrix} = \begin{bmatrix} \dot{x}_{k-1}^1 \\ \dot{x}_{k-1}^2 \end{bmatrix} + \mathbf{q}_{k-1}^l$$

$$\begin{bmatrix} x_k^1 \\ x_k^2 \end{bmatrix} = \begin{bmatrix} x_{k-1}^1 \\ x_{k-1}^2 \end{bmatrix} + T \begin{bmatrix} \dot{x}_{k-1}^1 \\ \dot{x}_{k-1}^2 \end{bmatrix} + \mathbf{q}_{k-1}^u$$

$$\mathbf{y}_k = \text{atan}_2(x_k^2, x_k^1) + \mathbf{r}_k$$

$$\text{where } \mathbf{r}_k \sim \mathcal{N}(0, (\frac{\pi}{180})^2) \text{ \& } \mathbf{q}_k = \begin{bmatrix} \mathbf{q}_k^u \\ \mathbf{q}_k^l \end{bmatrix} \sim \mathcal{N}\left(\mathbf{0}, \begin{bmatrix} \frac{T^2}{2} \mathbf{I} \\ \frac{T}{\Pi} \end{bmatrix} \begin{bmatrix} \underbrace{1}_{\text{red}} & 0 \\ 0 & \underbrace{1}_{\text{red}} \end{bmatrix} \begin{bmatrix} \frac{T^2}{2} \mathbf{I} \\ \frac{T}{\Pi} \end{bmatrix}^T\right)$$

Concluding remarks:

- Rao-Blackwellized particle filters are useful to reduce the number of particles.
- These filters enable us to handle higher dimensions than normal PFs.
- They are particularly useful if Kalman gains and posterior covariances are independent of the nonlinear states
⇒ sufficient to compute them one time in each recursion.