# CSC 445 - Intro to Intelligent Robotics, Spring 2018

Uncertainty

# **Probabilistic Robotics**

#### Probabilistic robotics is about the

- representation
- propagation
- reduction

of uncertainty.

# Environmental Representation

- The environment is characterized by *state*.
- There are two fundamental types of interactions between a robot and its environment.
  - Environmental sensor measurements
  - Control actions

# Notation

- The state is represented as  $x_t$  where t is time.
- The measurement data at time t is denoted as  $z_t$ .
- The control data at time t is denoted as  $u_t$ .
- The notation

$$y_{t_1:t_2} = y_{t_1}, y_{t_1+1}, \ldots, y_{t_2}$$

denotes the set of all y values between from  $t_1$  to  $t_2$ .

# Representing Uncertainty

- In probabilistic robotics, uncertainty is represented explicitly using probability theory.
- The evolution of a state may be represented as a probability distribution

$$p(x_t \mid x_{0:t-1}, z_{1:t-1}, u_{1:t})$$

That is, the state x<sub>t</sub> is conditioned on all past states, measurements and controls.

# Representing Uncertainty

The state transition probability

$$p(x_t \mid x_{0:t-1}, u_{1:t})$$

specifies how the state evolves over time as a function of the control actions.

The measurement probability

 $p(z_t \mid x_{0:t}, z_{1:t-1}, u_{1:t})$ 

specifies the probabilistic law according to which measurement z are generated from environment state x.

# State Evolution Bayes Network



# Complete State

- A state  $x_t$  is complete if it is the best predictor of the future.
- In other words, completeness means that knowledge of the past states, measurements, or controls carry no additional information that would help us predict the future.
- If x<sub>t</sub> is complete, the evolution of a state may be represented as a state transition probability

 $p(x_t \mid x_{t-1}, u_t)$ 

 Additionally, the measurement probability can be represented as

$$p(z_t \mid x_t)$$

# Belief Distribution

- A belief reflects the robot's internal knowledge about the state of the environment.
- Probabilistic robotics represents beliefs with conditional probability distributions.
- A belief distribution is a posterior probability over state variables conditioned on the available data

 $bel(x_t) = p(x_t | z_{1:t}, u_{1:t})$ 

# Bayes Filter Algorithm

function BAYES FILTER(
$$bel(x_{t-1}), u_t, z_t$$
)  
for all  $x_t$  do  
 $\overline{bel}(x_t) = \int p(x_t \mid u_t, x_{t-1})bel(x_{t-1})dx_{t-1}$   
 $bel(x_t) = \eta p(z_t \mid x_t)\overline{bel}(x_t)$   
return  $bel(x_t)$ 

# **Different Realizations**

- The Bayes filter is a framework for recursive state estimation.
- There are different realizations.
- Different properties
  - Linear vs. non-linear models for state transition and measurements.
  - Parametric vs. non-parametric
  - ...

# Gaussian Filters

 A Gaussian filter represents beliefs by multivariate normal distributions

$$p(x) = \det(2\pi\Sigma)^{\frac{1}{2}} \exp\{-\frac{1}{2}(x-\mu)^T\Sigma^{-1}(x-\mu)\}$$

The Gaussian distribution is unimodal, so the posterior represents a single hypothesis.

#### The Gaussian Distribution

#### A 1D Gaussian distribution is defined as

$$p(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{\frac{(x-\mu)^2}{2\sigma^2}}$$

• A n dimensional Gaussian distribution is defined as

$$p(x) = rac{1}{\sqrt{(2\pi)^n |\Sigma|}} e^{rac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)}$$

where x is a vector and  $\Sigma$  is a covariance matrix.

#### Covariance Matrix

When X is a vector, the variance is expressed as a covariance matrix ∑ where

$$\sigma_{ij} = \mathrm{E}[(x_i - \mu_i)(x_j - \mu_j)]$$

A covariance matrix has the form

$$\Sigma = \begin{bmatrix} \sigma_1^2 & \rho_{12}\sigma_1\sigma_2 & \dots & \rho_{1n}\sigma_1\sigma_n \\ \rho_{12}\sigma_1\sigma_2 & \sigma_2^2 & \dots & \rho_{2n}\sigma_2\sigma_n \\ \vdots & \vdots & \ddots & \vdots \\ \rho_{1n}\sigma_1\sigma_n & \rho_{2n}\sigma_2\sigma_n & \dots & \sigma_n^2 \end{bmatrix}$$

where  $\rho_{ij}$  corresponds to the degree of correlation between the variables  $X_i$  and  $X_j$ .

#### Properties of the Gaussian Distribution

Given two independent random variables,  $X \sim \mathcal{N}(\mu_x, \sigma_x^2)$  and  $Y \sim \mathcal{N}(\mu_y, \sigma_y^2)$ , then

$$aX + b \sim \mathcal{N}(a\mu_x + b, a^2\sigma_x^2)$$

and

$$Z = X + Y \sim \mathcal{N}(\mu_x + \mu_y, \sigma_x^2 + \sigma_y^2)$$

# Properties of the Gaussian Distribution



# Linear Approximation of a Nonlinear Function

- Given Y = f(X) with X and Y assumed to be Gaussian and  $f(\cdot)$  a nonlinear function
- Approximate the function with a first order Taylor series expansion

$$Y \approx f(\mu_x) + \frac{\partial f}{\partial X}\Big|_{x=\mu_x}(X-\mu_x)$$

# Linear Approximation of a Nonlinear Function



# Transforming Uncertainty

- Propagation of uncertainty is the effect of uncertainty of a random variable to the uncertainty of a function based on the random variable.
- Given a function

y = f(x)

that maps a random variable x, to a random variable y.

- Let the standard deviation of x be given by  $\sigma_x$ .
- We can calculate the variance of  $\sigma_{y}^{2}$  as

$$\sigma_y^2 = \left(\frac{\partial f}{\partial x}\right)^2 \sigma_x^2$$

# Transforming Uncertainty

- If the function is a multivariable function that maps n inputs to m outputs, then the variances become covariance matrices.
- The covariance matrix of of *y* can be calculated as

$$\Sigma_y = J \Sigma_x J^T$$

where J is an  $m \times n$  Jacobian matrix.

#### Jacobian Matrix

• Let f(x) be a vector-valued function

$$f(x) = \begin{bmatrix} f_1(x) \\ f_2(x) \end{bmatrix}$$

 Let the gradient operator be the vector of (first order) partial derivatives

$$\nabla_{\mathsf{X}} = \begin{bmatrix} \frac{\partial}{\partial_{\mathsf{X}_1}} & \frac{\partial}{\partial_{\mathsf{X}_2}} & \dots & \frac{\partial}{\partial_{\mathsf{X}_n}} \end{bmatrix}^{\mathsf{T}}$$

Then, the Jacobian matrix is defined as

$$F_{x} = \begin{bmatrix} f_{1}(x) \\ f_{2}(x) \end{bmatrix} \cdot \nabla_{x} = \begin{bmatrix} \frac{\partial f_{1}}{\partial x_{1}} \cdots \frac{\partial f_{1}}{\partial x_{n}} \\ \frac{\partial f_{2}}{\partial x_{1}} \cdots \frac{\partial f_{2}}{\partial x_{n}} \end{bmatrix}$$

# Kalman Filter

- The Kalman filter is a realization of a Bayes filter.
- It is an estimator for the linear Gaussian case.
- It is the optimal solution for linear models and Gaussian distributions.

#### Linear State Transition Model

■ The Kalman filter assumes that the true state as time k is evolved from the state at (k - 1) according to

$$x_k = F_k x_{k-1} + B_k u_k + w_k$$

where

- $F_k$  is the state transition model.
- $\blacksquare$   $B_k$  is the control-input model.
- *w<sub>k</sub>* is process noise assumed to be drawn from a zero mean normal distribution with covariance *Q<sub>k</sub>*.

# Linear Observation Model

At time k an observation (or measurement) zk of the true state xk is made according to

 $z_k = H_k x_k + v_k$ 

where

- *H<sub>k</sub>* is the observation model which maps the true state space into the observed state space.
- v<sub>k</sub> is observation noise assumed to be drawn from a zero mean normal distribution with covariance R<sub>k</sub>.

# Kalman Filter State

- The state of a Kalman filter is represented by two variables
  - \$\hat{x}\_{k|k}\$, the posterior state estimate at time k given the observations up to and including time k;
  - $P_{k|k}$ , the posterior error covariance matrix.
- The notation  $\hat{x}_{n|m}$  represents the estimate of x at time n given observations up to and including time  $m \le n$ .

#### Kalman Filter Prediction Step

#### 1 Predict the state estimate

$$\hat{x}_{k|k-1} = F_k \hat{x}_{k-1|k-1} + B_k u_k$$

2 Predict the estimate covariance

$$P_{k|k-1} = F_k P_{k-1|k-1} F_k^T + Q_k$$

# Kalman Filter Correction Step

**1** Compute the innovation

 $\tilde{y}_k = z_k - H_k \hat{x}_{k|k-1}$ 

2 Compute the innovation covariance

$$S_k = R_k + H_k P_{k|k-1} H_k^T$$

3 Compute the optimal Kalman gain

$$K_k = P_{k|k-1} H_k^T S_k^{-1}$$

4 Update the state estimate

$$\hat{x}_{k|k} = \hat{x}_{k|k-1} + K_k \tilde{y}_k$$

5 Update the estimate covariance

$$P_{k|k} = (I - K_k H_k) P_{k|k-1}$$

# The Extended Kalman Filter

- The Extended Kalman Filter (EKF) is a sub-optimal extension to the original Kalman filter algorithm.
- The EKF allows for the estimation of non-linear state transition and observation models.
- This is accomplished by linearizing the mean and covariance estimates.

#### Nonlinear State Transition Model

• The EKF assumes that the true state as time k is evolved from the state at (k - 1) according to

$$x_k = f(x_{k-1}, u_k) + w_k$$

where

- $f(\cdot)$  is the nonlinear state transition model.
- $F_k$  is the Jacobian of f with respect to the state.
- *w<sub>k</sub>* is process noise assumed to be drawn from a zero mean normal distribution with covariance *Q<sub>k</sub>*.

# Nonlinear Observation Model

At time k an observation (or measurement) zk of the true state xk is made according to

 $z_k = h(x_k) + v_k$ 

where

- h(·) is the nonlinear observation model which maps the true state space into the observed state space.
- $H_k$  is the Jacobian of h with respect to the state.
- v<sub>k</sub> is observation noise assumed to be drawn from a zero mean normal distribution with covariance R<sub>k</sub>.

# Extended Kalman Filter Prediction Step

$$\hat{x}_{k|k-1} = f(\hat{x}_{k-1|k-1}, u_k)$$

2 Predict the estimate covariance

$$P_{k|k-1} = F_k P_{k-1|k-1} F_k^T + Q_k$$

#### Extended Kalman Filter Correction Step

1 Compute the innovation

 $\tilde{y}_k = z_k - h(\hat{x}_{k|k-1})$ 

2 Compute the innovation covariance

$$S_k = H_k P_{k|k-1} H_k^T + R_k$$

3 Compute the (near optimal) Kalman gain

$$K_k = P_{k|k-1} H_k^T S_k^{-1}$$

4 Update the state estimate

$$\hat{x}_{k|k} = \hat{x}_{k|k-1} + K_k \tilde{y}_k$$

5 Update the estimate covariance

$$P_{k|k} = (I - K_k H_k) P_{k|k-1}$$