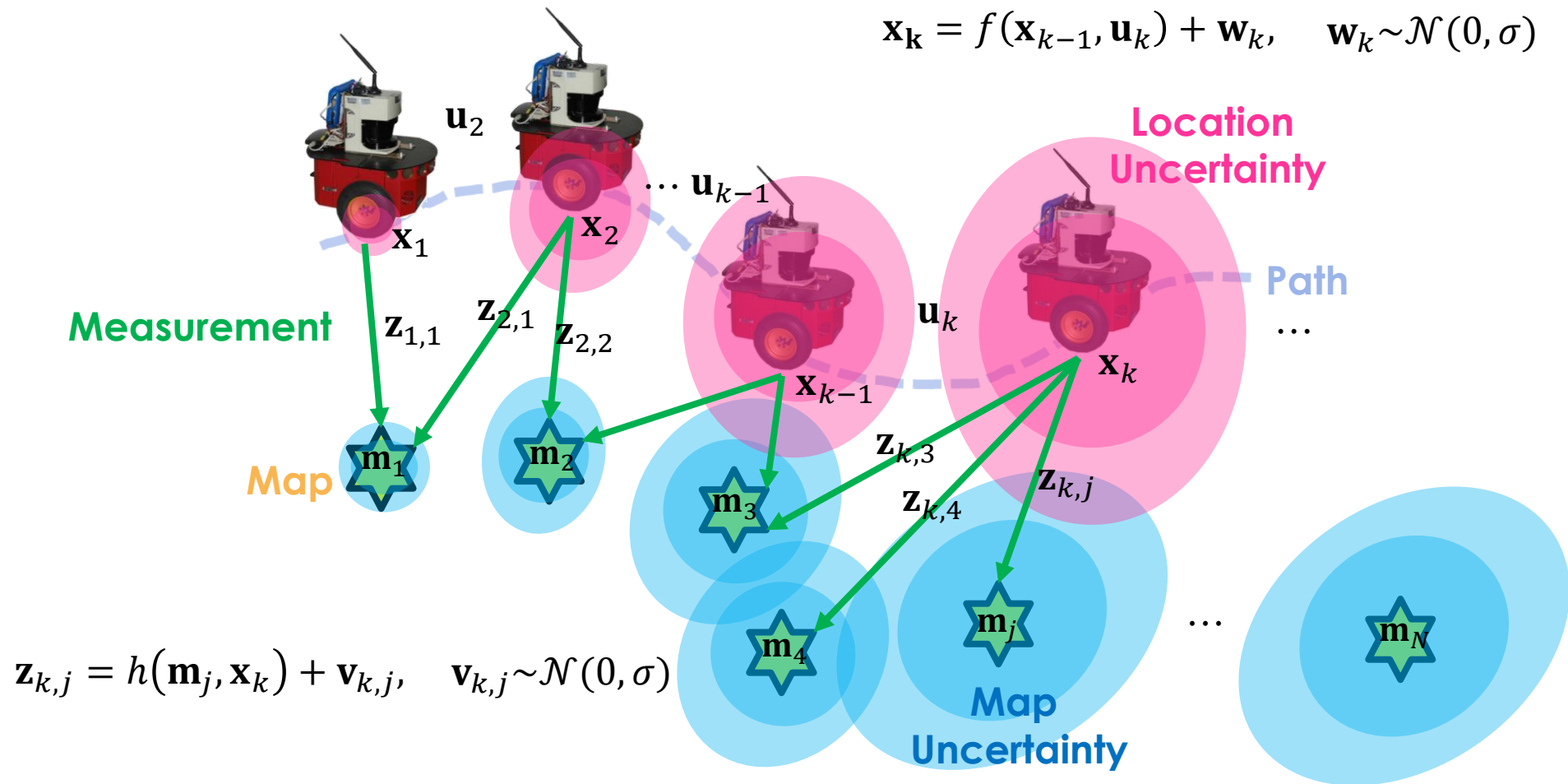


# Robotic Navigation and Exploration

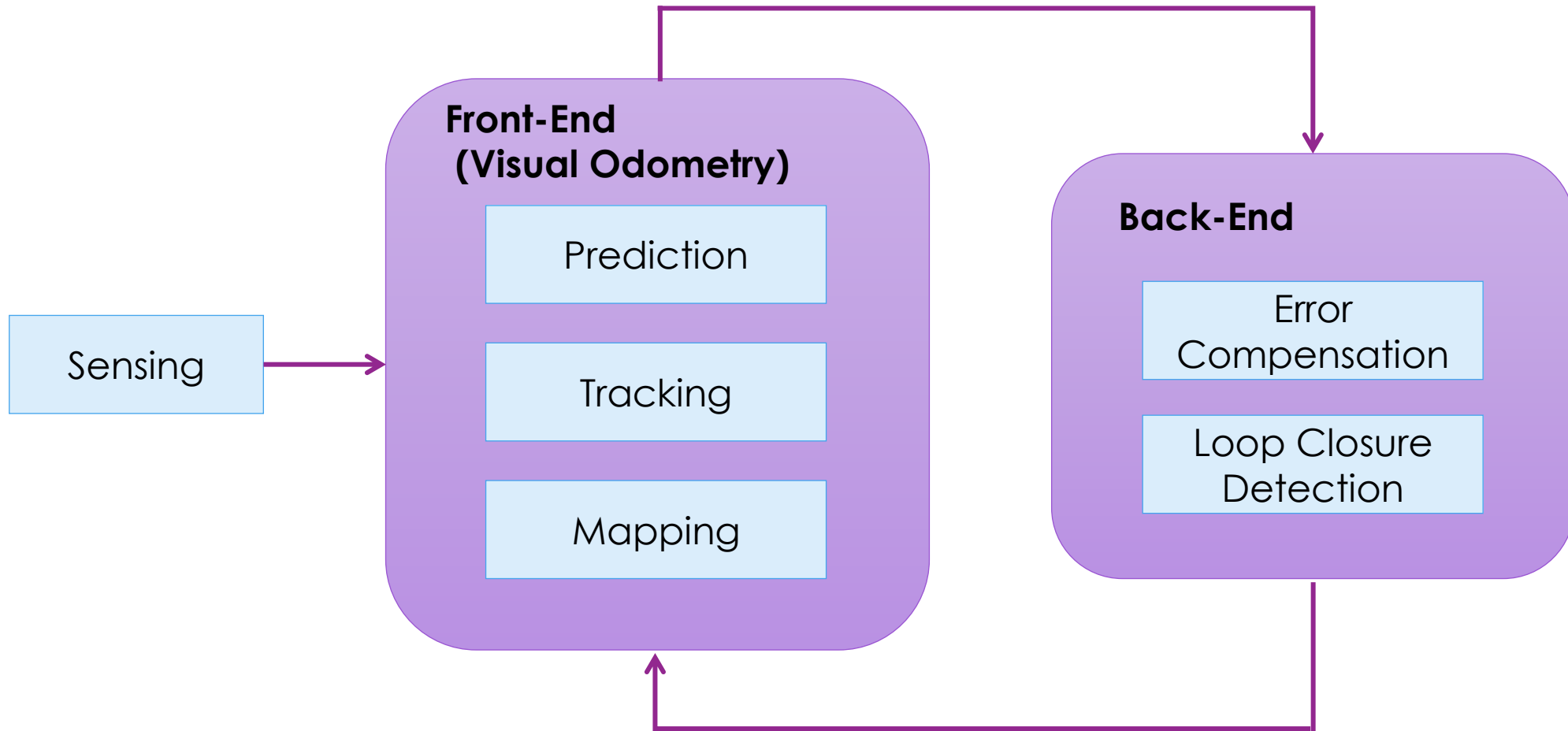
Week 5: SLAM Back-end (II)

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CS, NTHU

# Recall the SLAM Problem



# SLAM Architecture



# Error Compensation Methods

- Filter-based
  - Small Computation
  - On-line Optimization
- Graph-based
  - Large computation
  - High Accuracy
  - Off-line Optimization

# Outline

- State Estimation and SLAM Problem
- SLAM Back-end (Error Compensation)
  - Filter-based Methods
    - Probability Theory and Bayes Filter
    - Kalman Filter (KF) / Extended Kalman Filter (EKF)
      - EKF-SLAM
    - Particle Filter
      - Fast-SLAM
  - Graph-based Methods
    - Pose Graph and Least-square Optimization
    - Gauss-Newton and Levenberg-Marquardt Algorithm
    - Sparse Matrix for Optimization

# Introduction to Particle Filter

- EKF-SLAM assumes the probability distribution of robot pose and landmarks to be **Gaussian**, which leads to the following drawbacks:
  - Gaussian distribution can not express the robot pose properly.
  - The time complexity of estimating the covariance matrix for pose and landmarks is  $O(K^2)$ , which is time-consuming even when only observing few landmarks.  
( $K$ : number of landmarks)
- Particle filter utilizes **importance sampling** to approximate **arbitrary distribution**, which can express the robot pose more precisely.
- Furthermore, the time complexity of posterior estimation can be decreased to  **$O(M \log K)$**  by **disentangling the estimation process of pose and map**.

# Sampling Process

- In statistical modeling and inference, there are many complex problems that the closed-form descriptions of  $\mathbf{P}(\mathbf{X})$  can not be obtained.
- One can define a function  $f(X)$  that computes  $P(X)$  up to a normalizing constant:

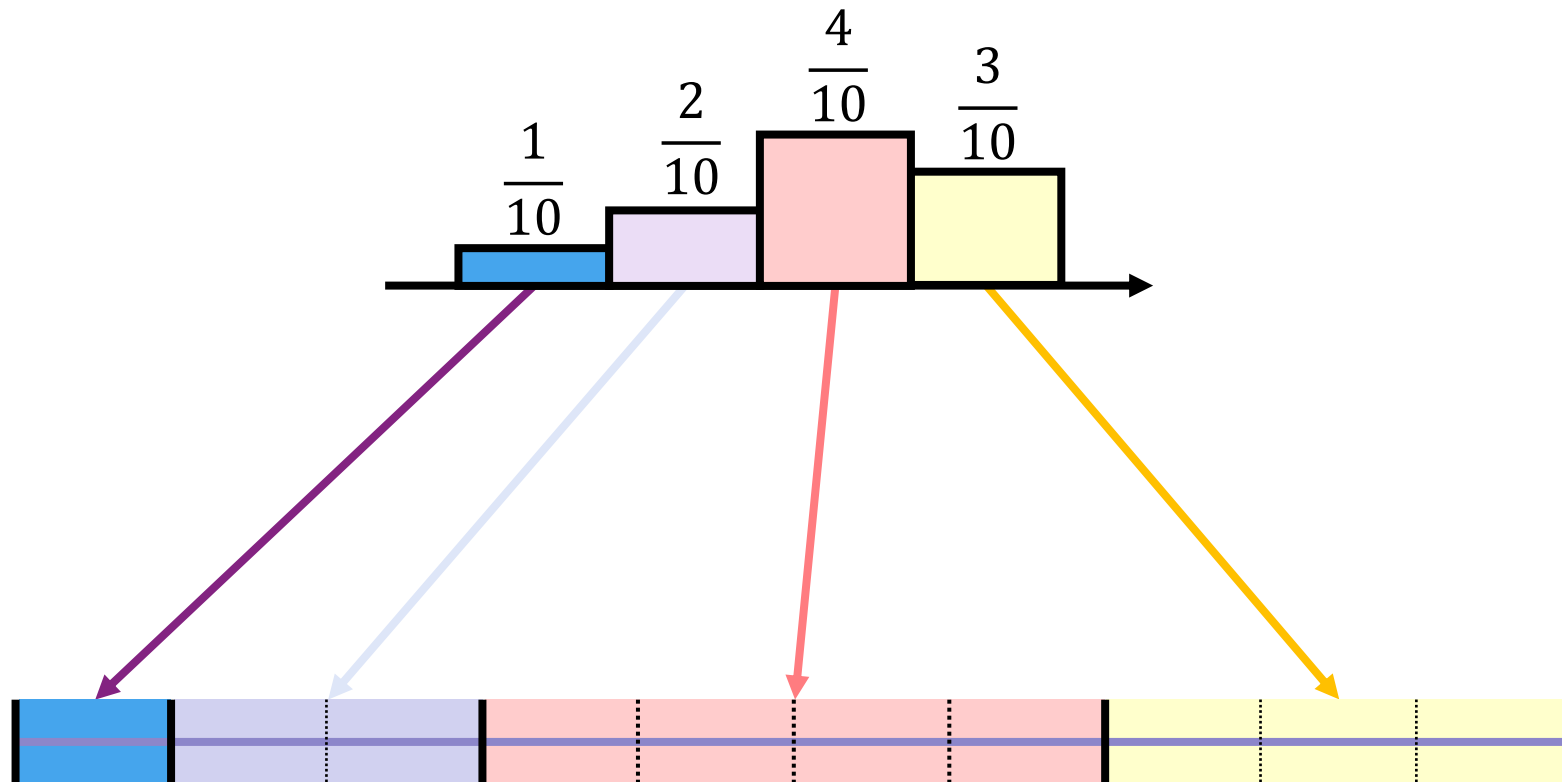
$$p(X) = \frac{f(X)}{Z}$$

where  $z = \int_{x \in \mathcal{S}} f(x) dx$  can not be computed because  $f(X)$  is too complex, or because the state space  $\mathcal{S}$  is too large to compute the integral.

- Statistical sampling and simulation techniques are used for getting fair samples from target probability distributions.

# Basic Sampling

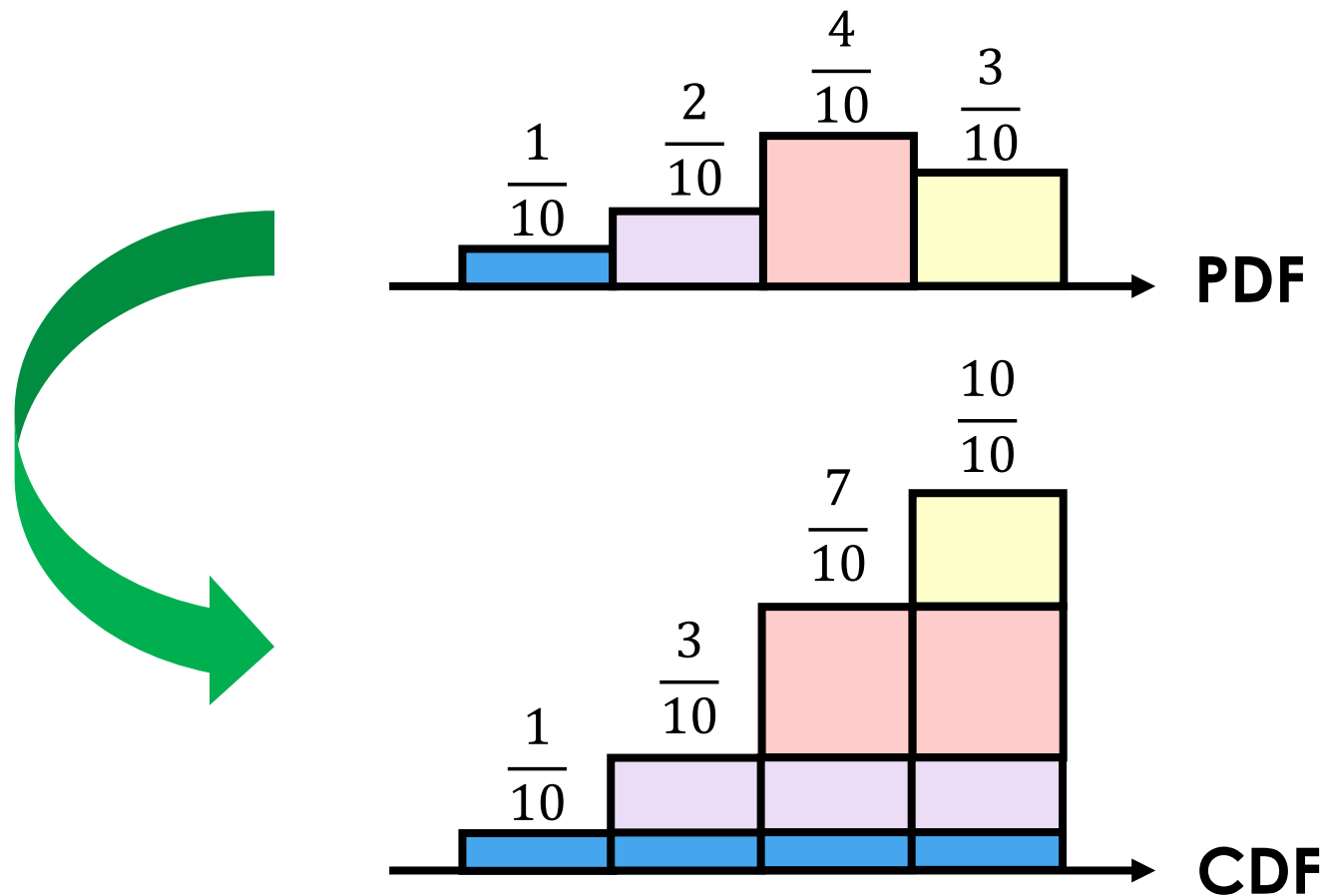
- Sampling from Probability Distribution Figure (PDF)





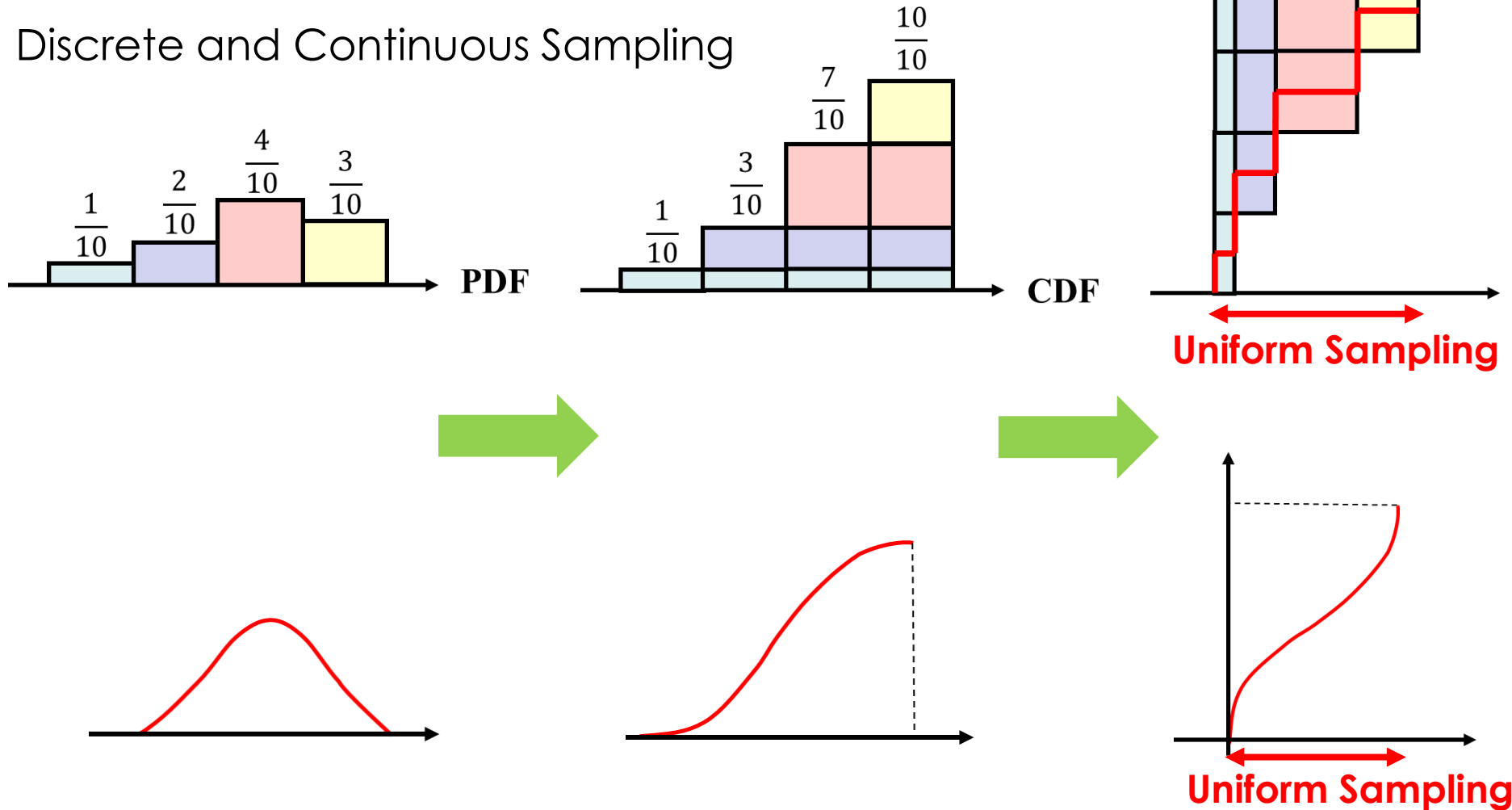
# Basic Sampling

- From Probability Distribution Figure (PDF) to Cumulated Distribution Figure (CDF)



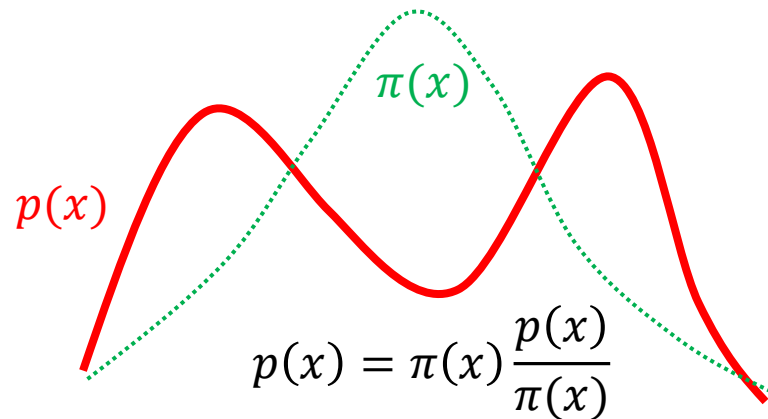
# Basic Sampling

- Discrete and Continuous Sampling

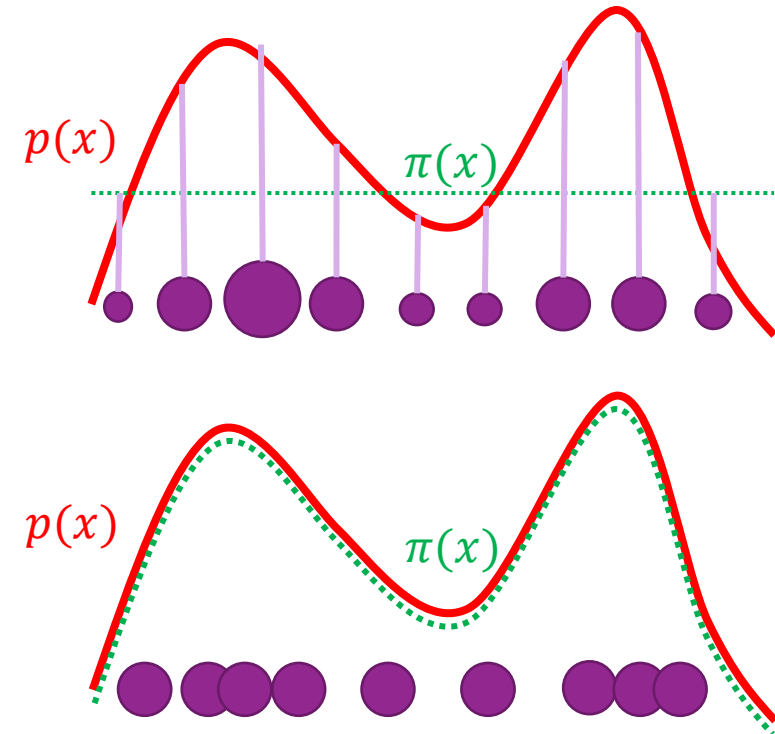


# Importance Sampling

- Important sampling adopts discrete multinomial to approximate arbitrary distribution. More sampling particles will have more accurate approximation.



1. Sampling  $x_i$  from  $\pi(x)$
2. Calculate  $w_i = \frac{p(x_i)}{\pi(x_i)}$
3. Sampling  $x$  from  $\text{mul}(x_i, w_i)$



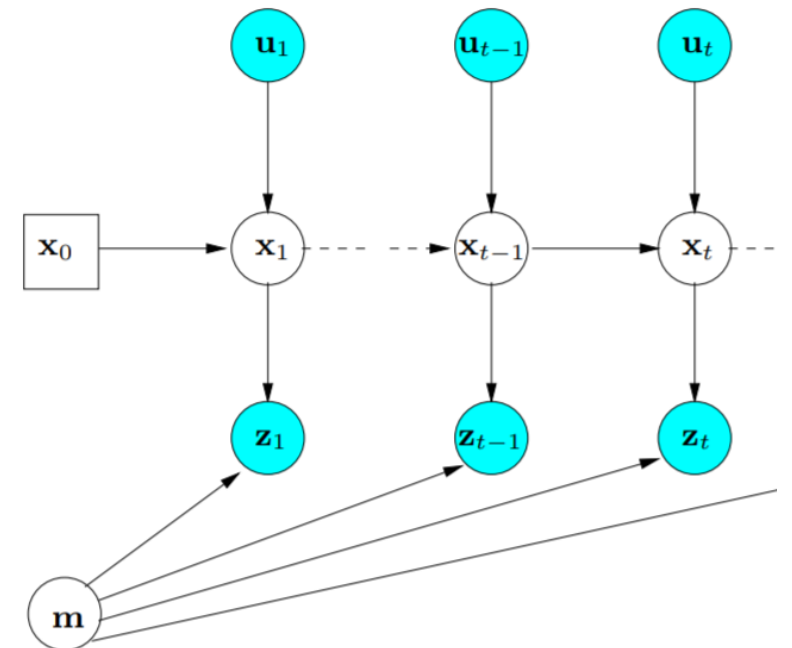
# Sequential Importance Sampling (SIS)

- Consider the localization problem, we utilize several particles to represent the approximation of pose distribution.
- In importance sampling, each particle have its own pose and weighting. The weighting is the division of source distribution and target distribution:

$$w_t^{(i)} = \frac{p(x_{1:t}^{(i)} | z_{1:t}, u_{1:t-1})}{\pi(x_{1:t}^{(i)} | z_{1:t}, u_{1:t-1})}$$

- According to the graphical model, we have

$$w_t^{(i)} = \frac{p(x_t^{(i)} | x_{1:t-1}^{(i)}, z_{1:t}, u_{1:t-1})}{\pi(x_t^{(i)} | x_{1:t-1}^{(i)}, z_{1:t}, u_{1:t-1})} \cdot \frac{p(x_{1:t-1}^{(i)} | z_{1:t-1}, u_{1:t-2})}{\pi(x_{1:t-1}^{(i)} | z_{1:t-1}, u_{1:t-2})}$$



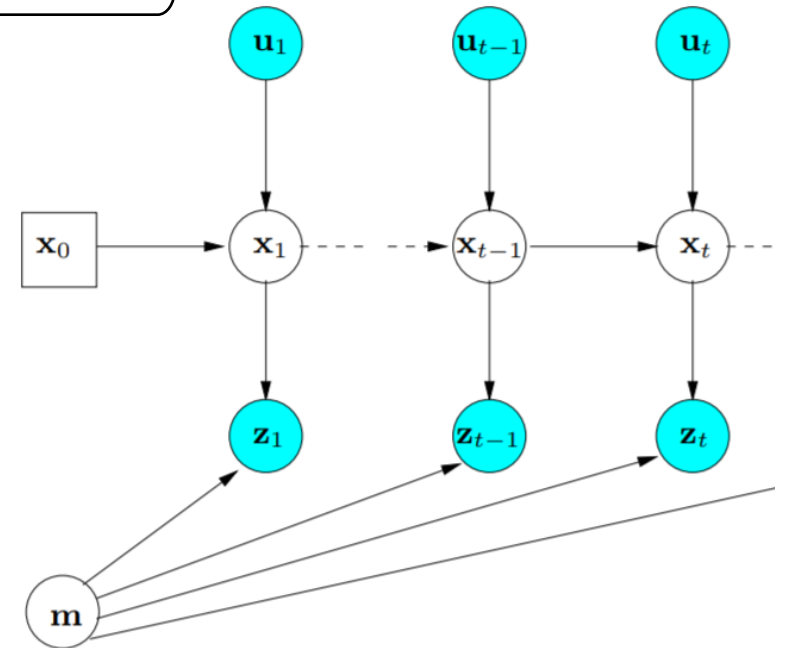
# Sequential Importance Sampling (SIS)

- Apply the Bayes theorem, we can get

$$w_t^{(i)} = \frac{\eta p(z_t | x_{1:t}^{(i)}, u_{1:t-1}) p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})}{\pi(x_t^{(i)} | x_{1:t-1}^{(i)}, z_{1:t}, u_{1:t-1})} \cdot \underbrace{\frac{p(x_{1:t-1}^{(i)} | z_{1:t-1}, u_{1:t-2})}{\pi(x_{1:t-1}^{(i)} | z_{1:t-1}, u_{1:t-2})}}_{w_{t-1}^{(i)}}$$

$$\propto \frac{p(z_t | m_{t-1}, x_t^{(i)}) p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})}{\pi(x_t^{(i)} | x_{1:t-1}^{(i)}, z_{1:t}, u_{1:t-1})} \cdot w_{t-1}^{(i)}$$

, in which  $\eta = \frac{1}{p(z_t | z_{1:t-1}, u_{1:t-1})}$



# Sequential Importance Sampling (SIS)

- Now, we select the distribution of last timestep as the source distribution:

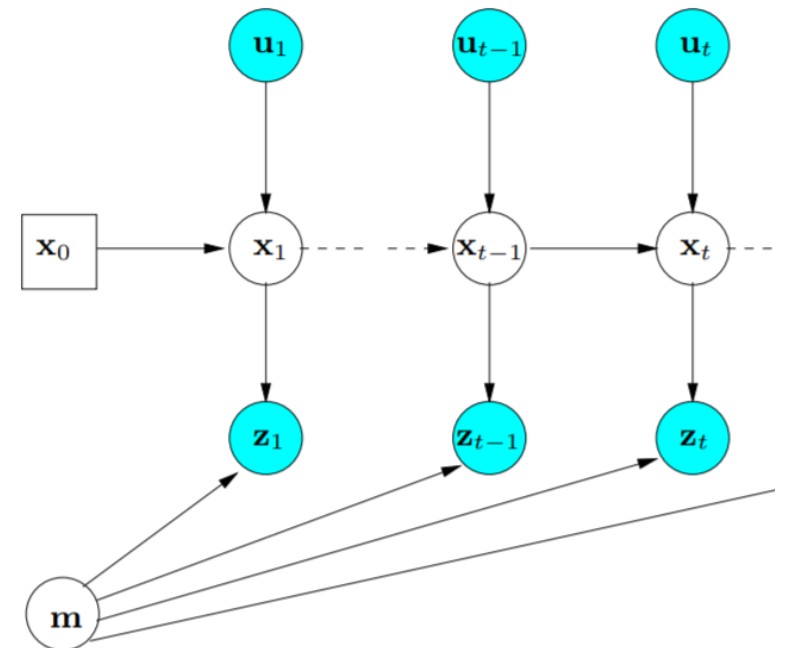
$$\pi(x_t | x_{1:t-1}, z_{1:t}, u_{1:t-1}) = p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})$$

$$w_t^{(i)} = \frac{\eta p(z_t | m_{t-1}, x_t^{(i)}) p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})}{\pi(x_t | x_{1:t-1}, z_{1:t}, u_{1:t-1})} \cdot w_{t-1}^{(i)}$$

- We can get the update weighting:

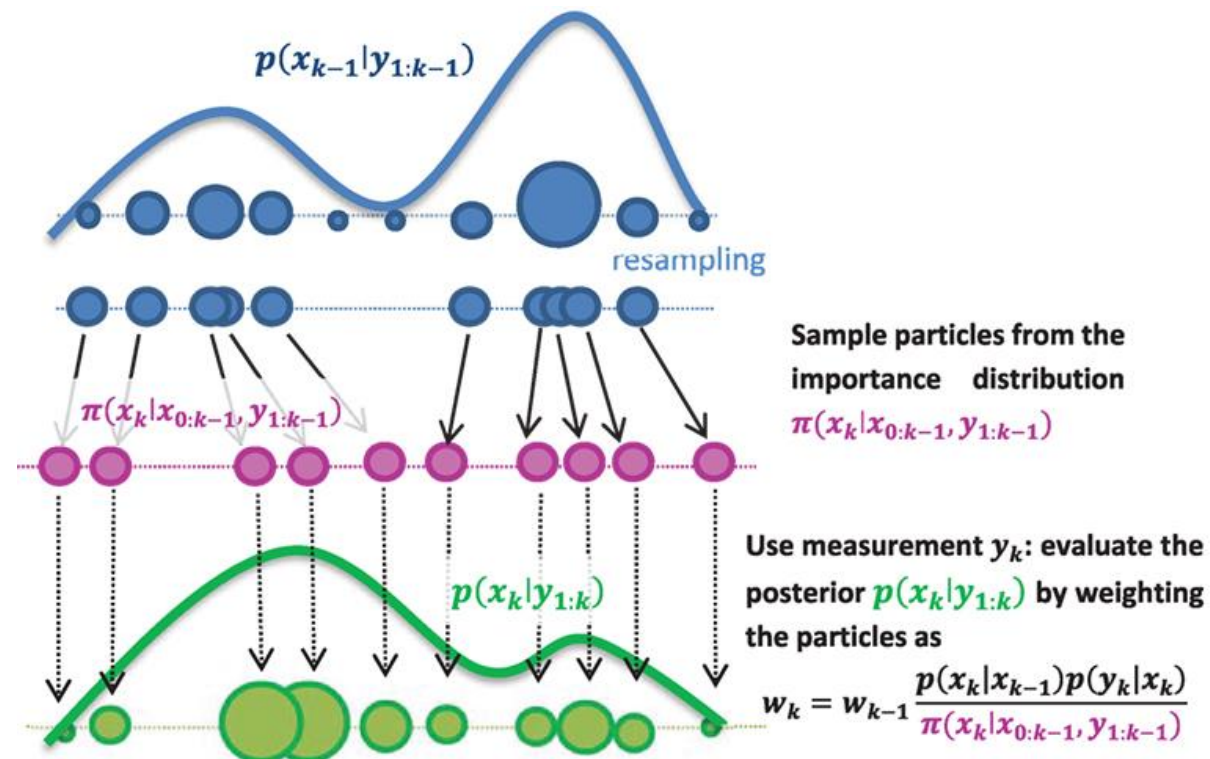
$$w_t^{(i)} = w_{t-1}^{(i)} \frac{\eta p(z_t | m_{t-1}, x_t^{(i)}) p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})}{p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})}$$

$$\propto w_{t-1}^{(i)} \cdot p(z_t | m_{t-1}, x_t^{(i)})$$



# Sequential Importance Resampling (SIR)

- After several steps, the weightings of most particles in SIS particle filter will decrease to close to zero.
- To avoid this problem, we can utilize the resampling process:



# Monte-Carlo Localization Example





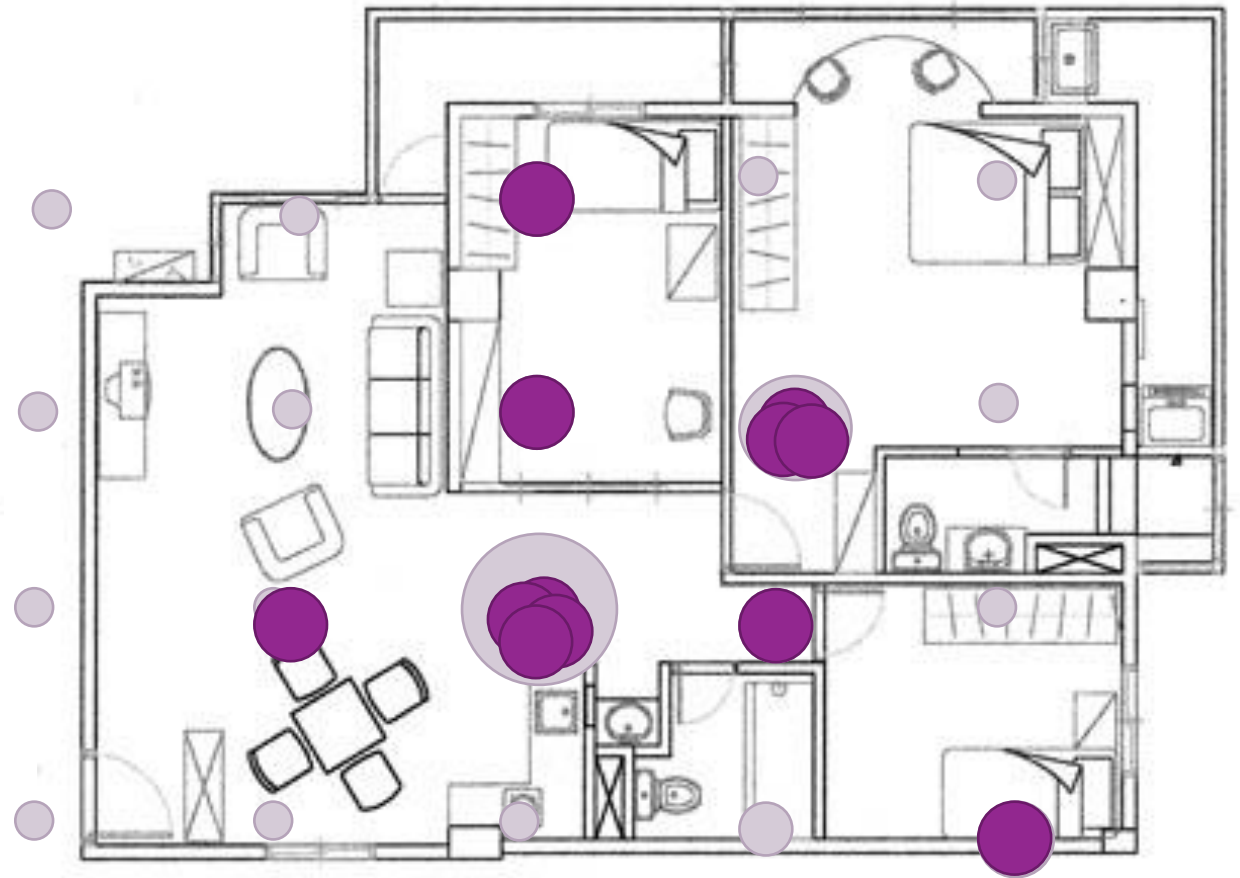
# Monte-Carlo Localization Example



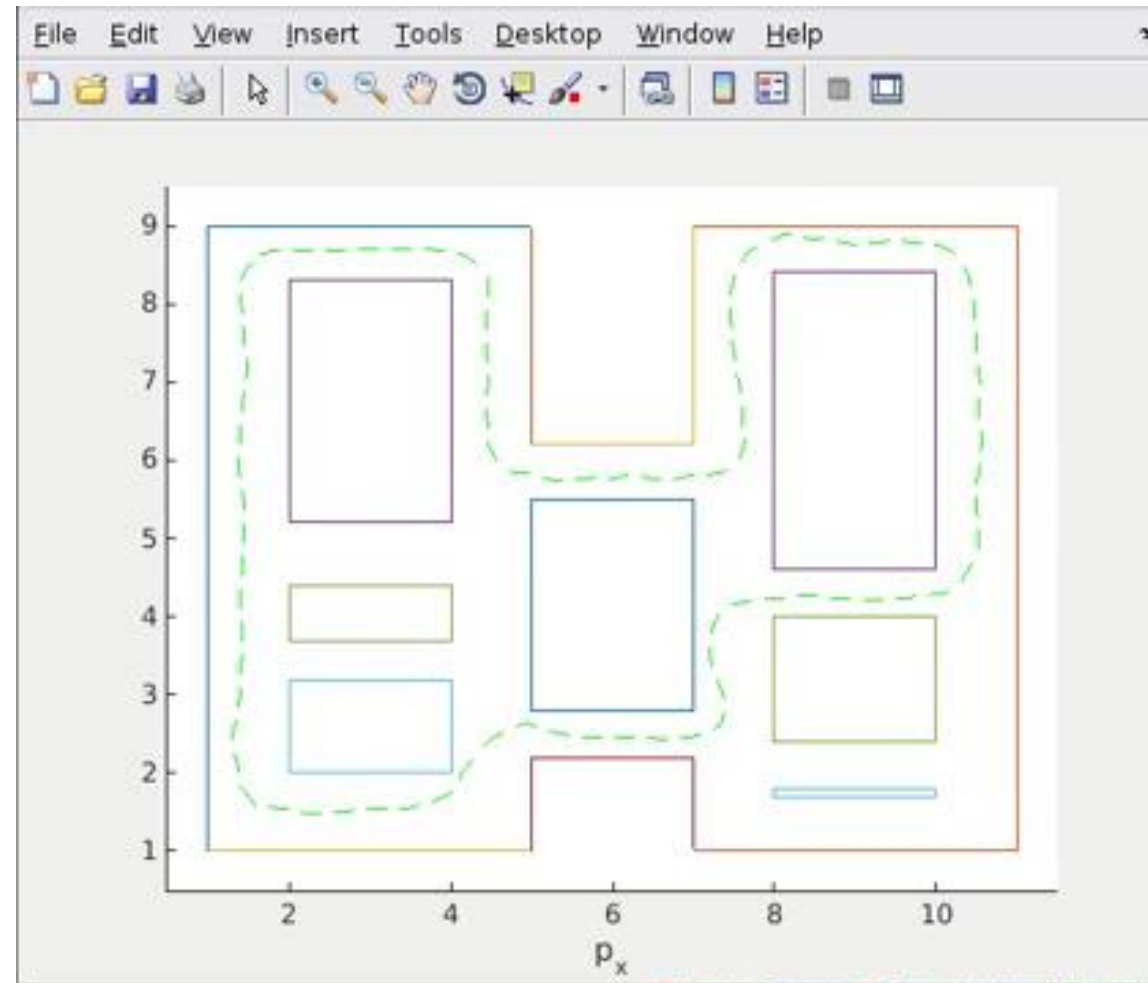
# Monte-Carlo Localization Example



# Monte-Carlo Localization Example



# Monte-Carlo Localization



# Fast-SLAM

- Now consider the full SLAM problem (localization and mapping), we can divide the full process to localization and mapping steps. This method is called **Rao-Blackwellization**.

$$p(x_{1:t}, m_t | z_{1:t}, u_{1:t}) = p(x_{1:t} | z_{1:t}, u_{1:t}) p(m_t | x_{1:t}, z_{1:t})$$

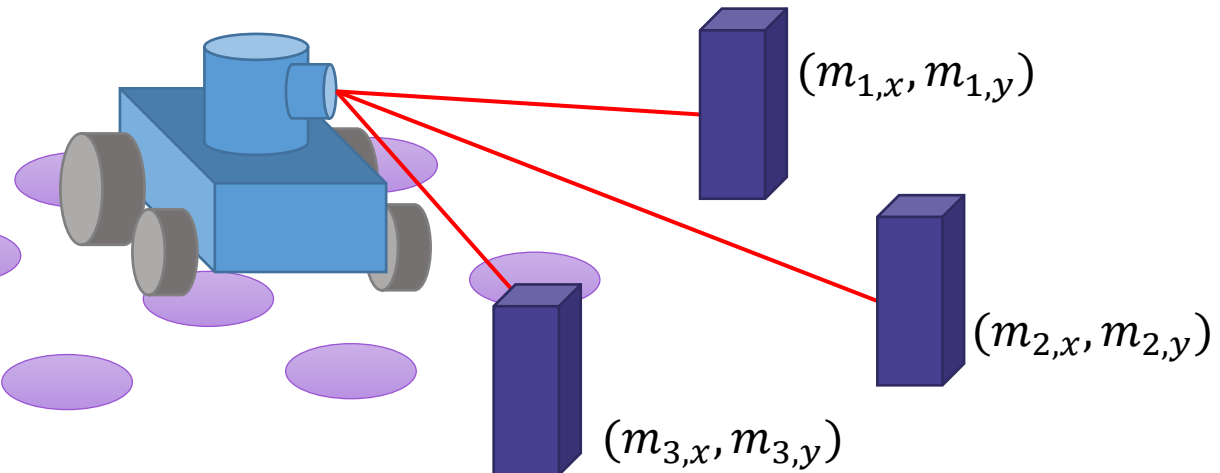
- In Fast-SLAM, the robot pose is represented by the multivariate distribution of several weighted particles, and each particle adopts **K** extended Kalman filter to estimate the landmarks independently.

Particle Weights:  $w^{(i)}$

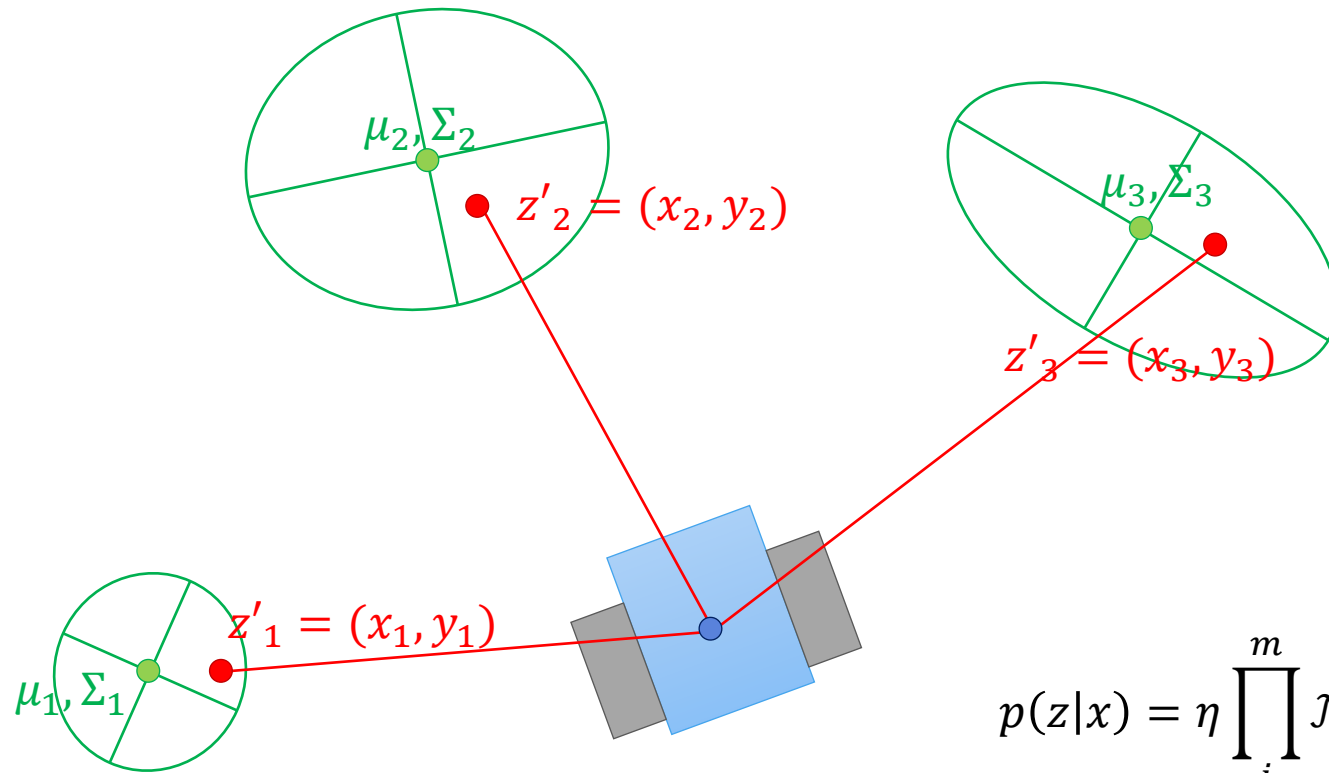
Robot Pose:  $(x \ y \ \theta)^{(i)}$

Landmarks:

$(\mu_1^{(i)}, \Sigma_1^{(i)}), (\mu_2^{(i)}, \Sigma_2^{(i)}), (\mu_3^{(i)}, \Sigma_3^{(i)})$



# Likelihood of Measurement



$$p(z|x) = \eta \prod_i^m \mathcal{N}(z'_i; \mu_i, \Sigma_i)$$

# Fast-SLAM

- Steps of Fast-SLAM

1. Predict the next pose  $x_t^{(i)}$  by motion model.

$$x_t^{(i)} \sim p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})$$

2. Update the distribution of each landmark  $(\mu_{j,t}^{(i)}, \Sigma_{j,t}^{(i)})$  via measurement  $z_k$ .

$$Q = H \Sigma_{j,t-1}^{(i)} H^T + R, \quad K_t = \Sigma_{j,t-1}^{(i)} H^T Q^{-1}$$

$$\mu_{j,t}^{(i)} = \mu_{j,t-1}^{(i)} + K_k \left( z_k - h(\mu_{j,t-1}^{(i)}, x_t^{(i)}) \right)$$

$$\Sigma_{j,t}^{(i)} = (I - K_t H) \Sigma_{j,t-1}^{(i)}$$

3. Update the importance weight of particles.

$$w^{(i)} \sim |2\pi Q|^{-\frac{1}{2}} \exp \left\{ -\frac{1}{2} \left( z_k - h(\mu_{j,t-1}^{(i)}, x_t^{(i)}) \right)^T Q^{-1} \left( z_k - h(\mu_{j,t-1}^{(i)}, x_t^{(i)}) \right) \right\}$$

4. Resampling.



# Fast SLAM

- Measure of how well the target distribution is approximated by samples drawn from the proposal.

$$N_{eff} = \frac{1}{\sum_i (w_t^{(i)})^2}$$

- $N_{eff}$  denotes the inverse variance of the normalized particle weights. For equal weights, the results is the number of the particles. And the sample approximation is close to the target.

$$N_{eff}^* = \frac{1}{\sum_i \frac{1}{N^2}} = \frac{1}{N \frac{1}{N^2}} = N$$

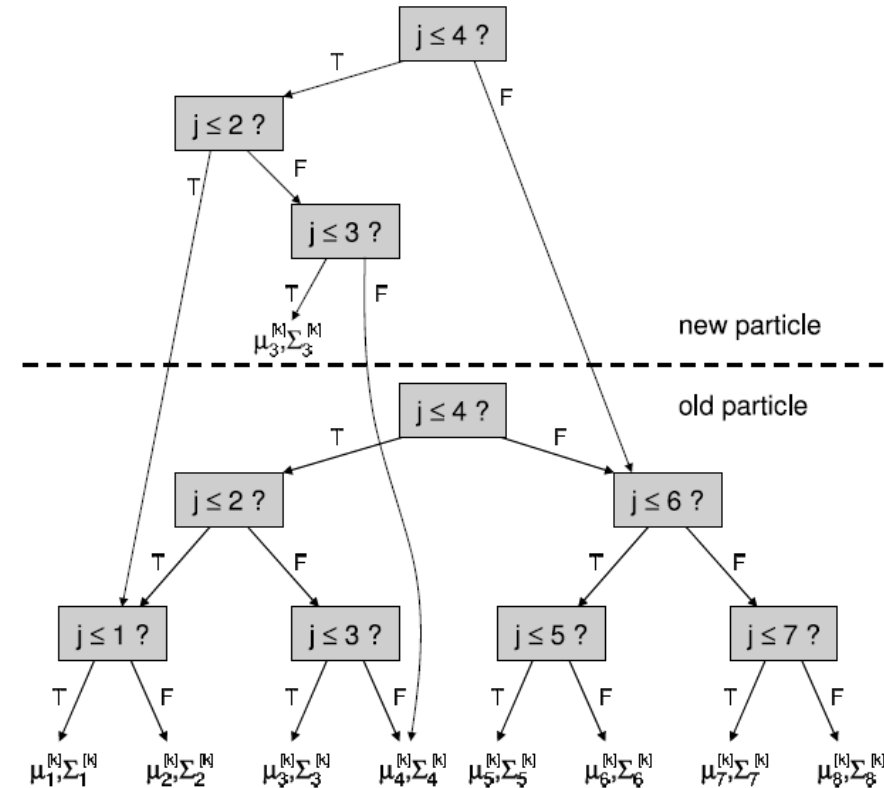
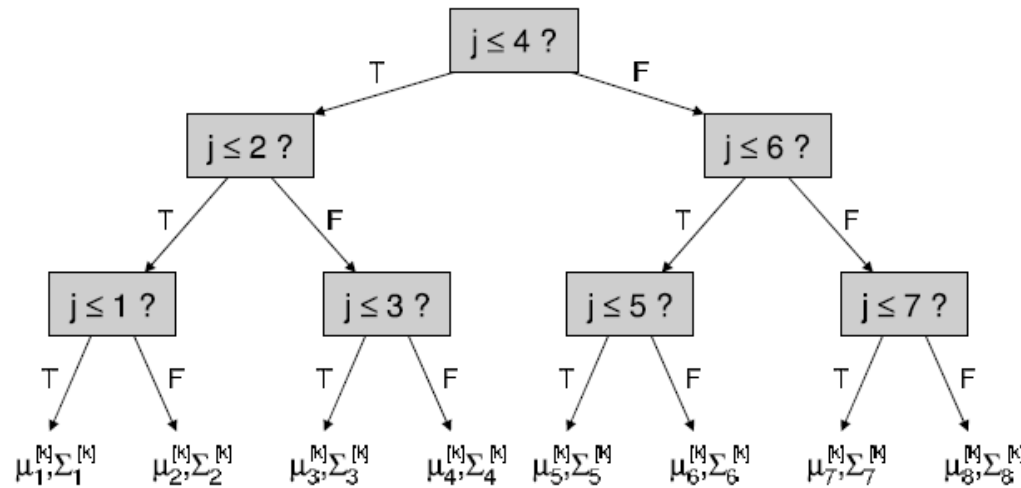
- If  $N_{eff}$  drops below a given threshold (usually set to half of the particles), we will resample the particle.

$$N_{eff} < \frac{N}{2}$$

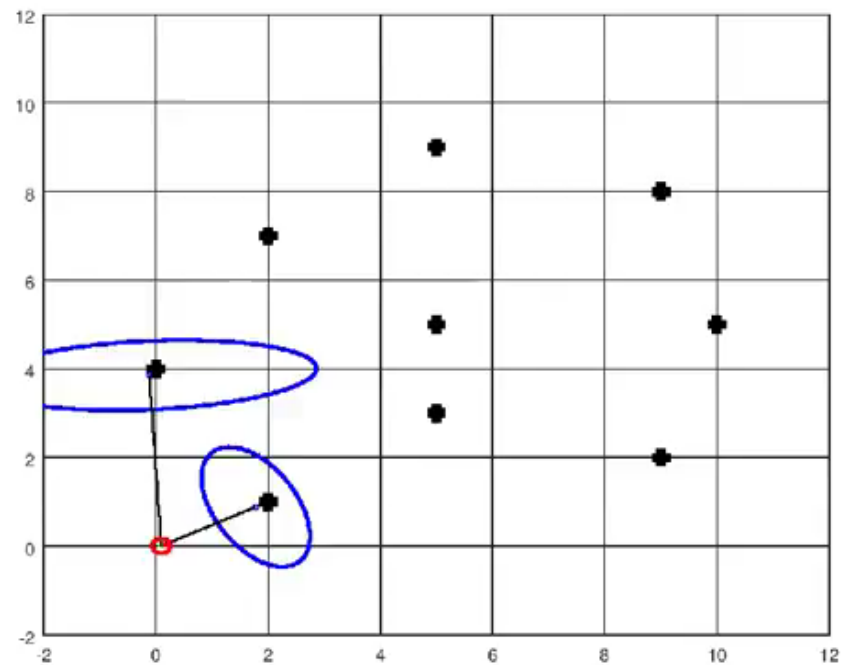


# Fast-SLAM

- Efficient implementation of Fast-SLAM. The basic idea is that the set of Gaussians in each particle is represented by a balanced binary tree.



# Fast-SLAM Demo



# Occupancy Grid Representation

- Occupancy grid maps is a volumetric representation of environments. The advantage of grid representation is that it does not require any predefined definition of landmarks. Instead, it can model arbitrary types of environments.



# Grid-based Fast-SLAM

- Steps of Grid-based Fast-SLAM

1. Predict the next pose  $x_t^{(i)}$  by motion model.
2. Update the occupancy grid map of each particle.
3. Update the importance weight of particles.
4. Resampling.

# Grid-based Fast-SLAM

- Steps of Grid-based Fast-SLAM

1. Predict the next pose  $x_t^{(i)}$  by motion model.
2. Update the occupancy grid map of each particle.
3. Update the importance weight of particles.
4. Resampling.

# Occupancy Grid Map Algorithm

- The occupancy grids store the probability if the discrete location is free. The state of the grid is defined by the rate of free and occupied.

$$Odd(s) = \frac{p(s = 1)}{p(s = 0)}$$

- Apply the Bayes theorem to compute the posterior of the state

$$p(s|z) = \frac{p(z|s)p(s)}{p(z)} \quad Odd(s|z) = \frac{p(s = 1|z)}{p(s = 0|z)} = \frac{p(z|s = 1)p(s = 1)/p(z)}{p(z|s = 0)p(s = 0)/p(z)} = \frac{p(z|s = 1)}{p(z|s = 0)} Odd(s)$$

- Utilize the log operation to simplify the computation

$$\log Odd(s|z) = \log \frac{p(z|s = 1)}{p(z|s = 0)} + \log Odd(s)$$

- Compute the occupied probability from log state

define  $g = \log Odd(s)$  and  $p = p(s = 1)$

$$\exp(g) = Odd(s) = \frac{p(s = 1)}{p(s = 0)} = \frac{p}{1 - p} \quad p = \frac{\exp(g)}{1 + \exp(g)}$$

# Occupancy Grid Map Algorithm

- Define two likelihood parameter

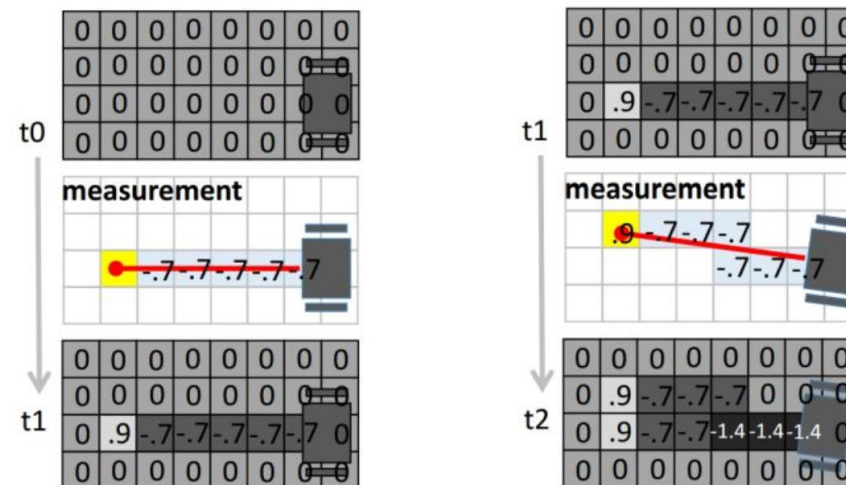
$$l_{occ} = \frac{p(z = 0 | s = 1)}{p(z = 0 | s = 0)} \quad l_{free} = \frac{p(z = 1 | s = 1)}{p(z = 1 | s = 0)}$$

$$\log Odd(s|z) = \log \frac{p(z|s = 1)}{p(z|s = 0)} + \log Odd(s)$$

- Initialize the state with half probability of occupied

$$\log Odd(s_{init}) = \log \frac{p(s_{init} = 1)}{p(s_{init} = 0)} = \log \frac{0.5}{0.5} = 0$$

- Ray tracing the grid, add  $l_{free}$  to all the grids passing by, and add  $l_{occ}$  to the last grid.



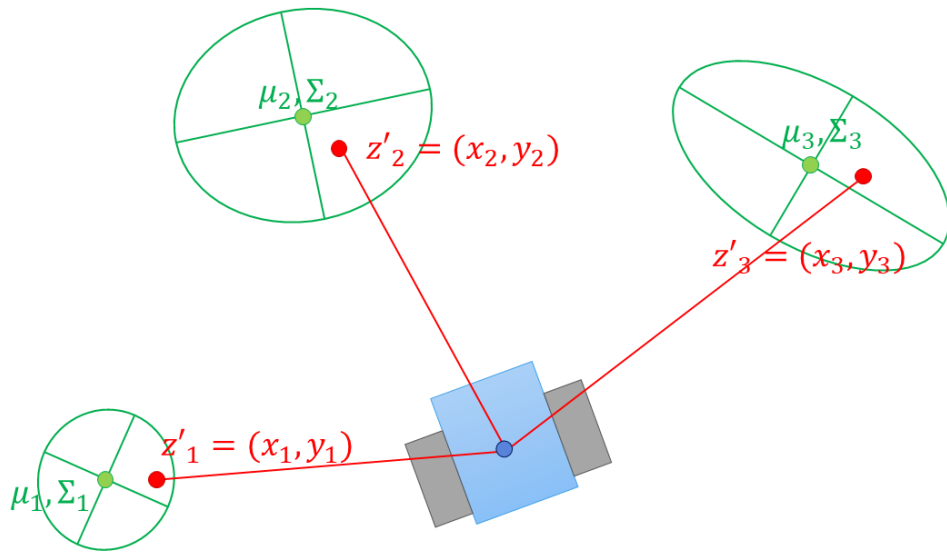
# Grid-based Fast-SLAM

- Steps of Grid-based Fast-SLAM

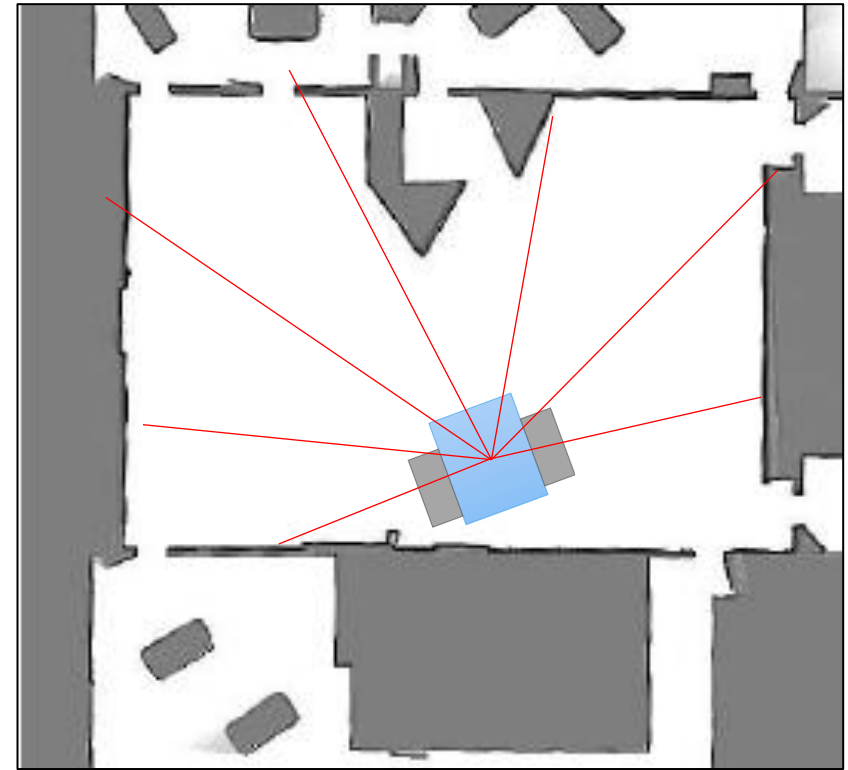
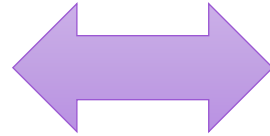
1. Predict the next pose  $x_t^{(i)}$  by motion model.
2. Update the occupancy grid map of each particle.
3. Update the importance weight of particles.
4. Resampling.



# Likelihood Field of Grid Map



$$p(z|x) = \eta \prod_i^m \mathcal{N}(z'_i; \mu_i, \Sigma_i)$$



$$p(z|x) = ?$$

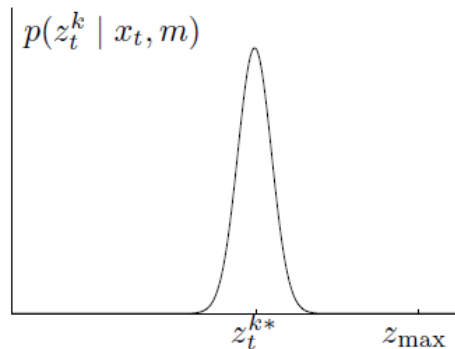
# Laser Beam Model

- A common sensor of mobile robot is a range finder, which measures the distance from the robot to the obstacles.

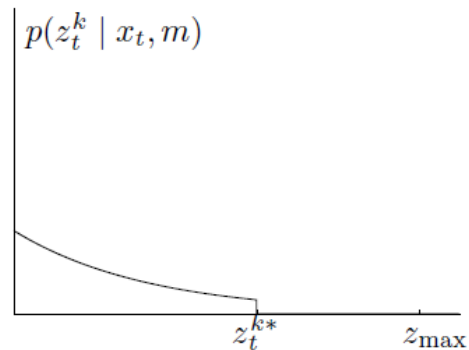


- 4 components of the measurement model:

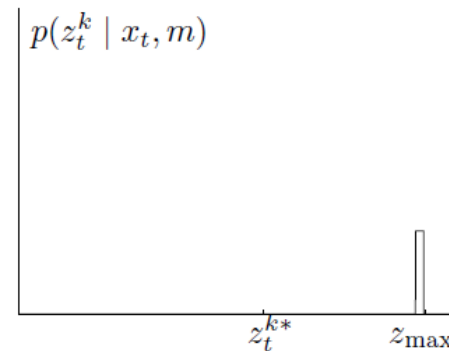
(a) Gaussian distribution  $p_{\text{hit}}$



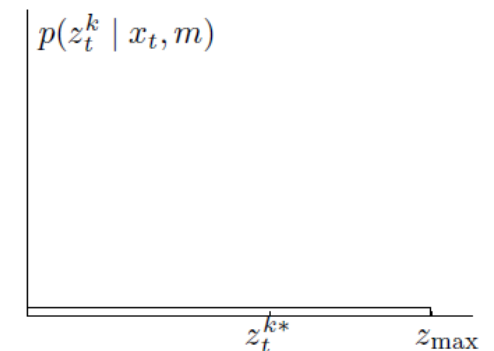
(b) Exponential distribution  $p_{\text{short}}$



(c) Uniform distribution  $p_{\text{max}}$



(d) Uniform distribution  $p_{\text{rand}}$



# Laser Beam Model

- (a) Correct range with local measurement noise.

$$\mathcal{N}(z_t^k; z_t^{k*}, \sigma_{\text{hit}}^2) = \frac{1}{\sqrt{2\pi\sigma_{\text{hit}}^2}} e^{-\frac{1}{2} \frac{(z_t^k - z_t^{k*})^2}{\sigma_{\text{hit}}^2}}$$

- (b) Unexpected objects

$$p_{\text{short}}(z_t^k | x_t, m) = \begin{cases} \eta \lambda_{\text{short}} e^{-\lambda_{\text{short}} z_t^k} & \text{if } 0 \leq z_t^k \leq z_t^{k*} \\ 0 & \text{otherwise} \end{cases}$$

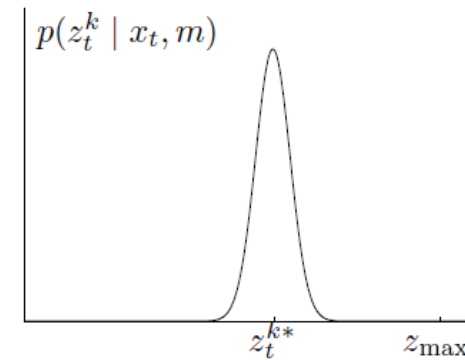
- (c) Failures

$$p_{\text{max}}(z_t^k | x_t, m) = I(z = z_{\text{max}}) = \begin{cases} 1 & \text{if } z = z_{\text{max}} \\ 0 & \text{otherwise} \end{cases}$$

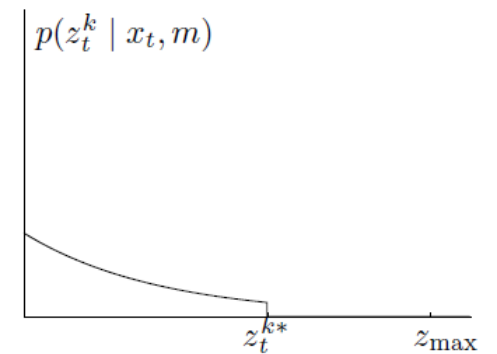
- (d) Random measurements

$$p_{\text{rand}}(z_t^k | x_t, m) = \begin{cases} \frac{1}{z_{\text{max}}} & \text{if } 0 \leq z_t^k < z_{\text{max}} \\ 0 & \text{otherwise} \end{cases}$$

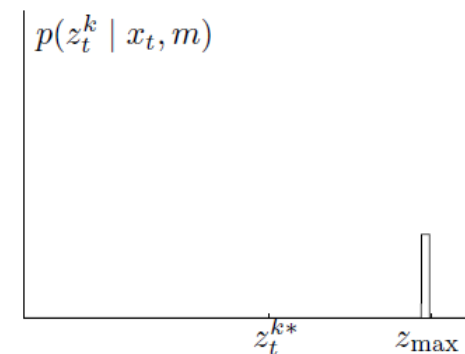
(a) Gaussian distribution  $p_{\text{hit}}$



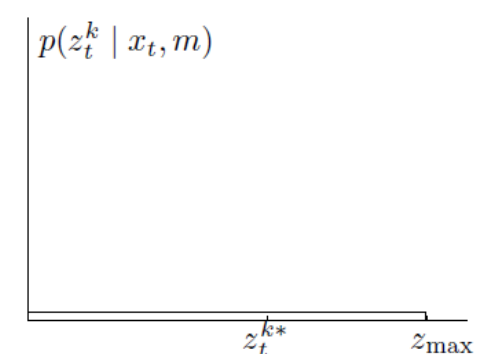
(b) Exponential distribution  $p_{\text{short}}$



(c) Uniform distribution  $p_{\text{max}}$

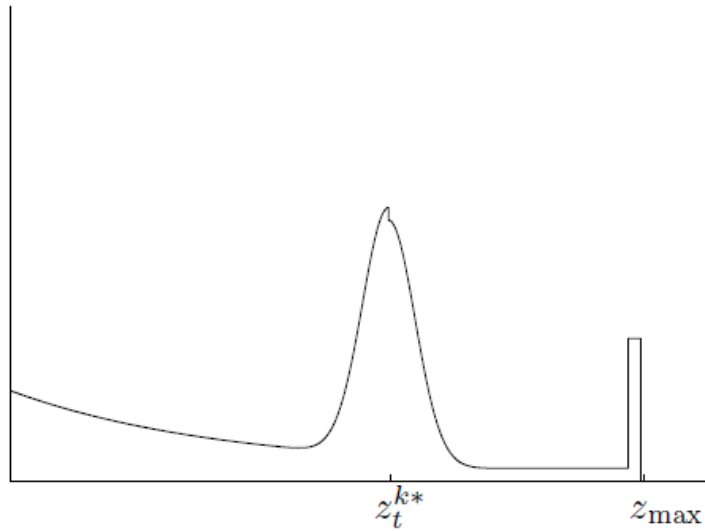


(d) Uniform distribution  $p_{\text{rand}}$



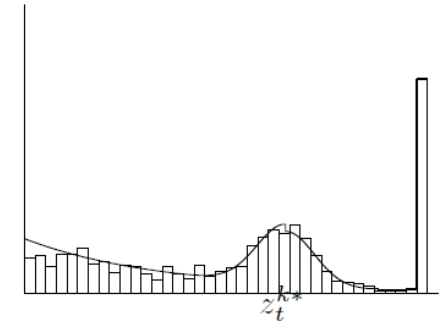
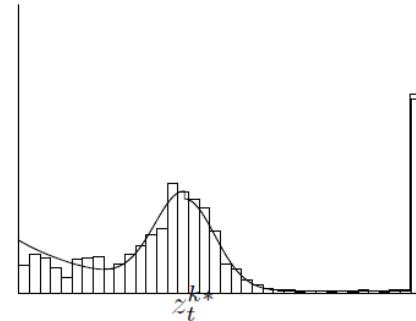
# Laser Beam Model

- Mixture distribution of laser beam model.

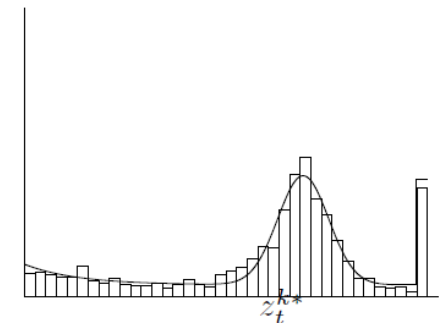
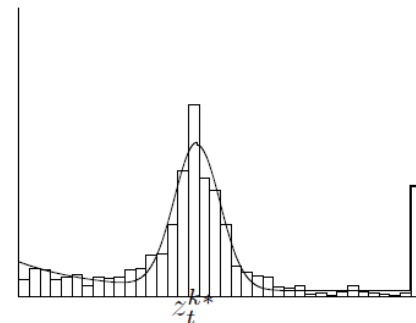


$$p(z_t^k | x_t, m) = \begin{pmatrix} z_{\text{hit}} \\ z_{\text{short}} \\ z_{\text{max}} \\ z_{\text{rand}} \end{pmatrix}^T \cdot \begin{pmatrix} p_{\text{hit}}(z_t^k | x_t, m) \\ p_{\text{short}}(z_t^k | x_t, m) \\ p_{\text{max}}(z_t^k | x_t, m) \\ p_{\text{rand}}(z_t^k | x_t, m) \end{pmatrix}$$

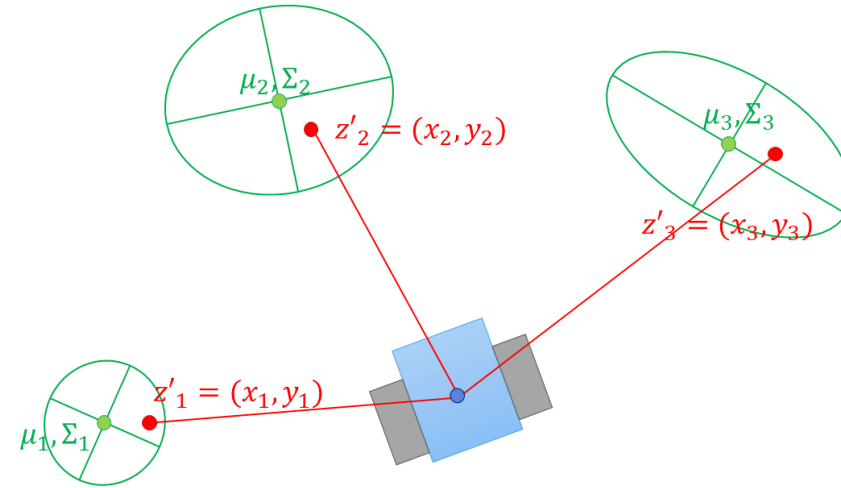
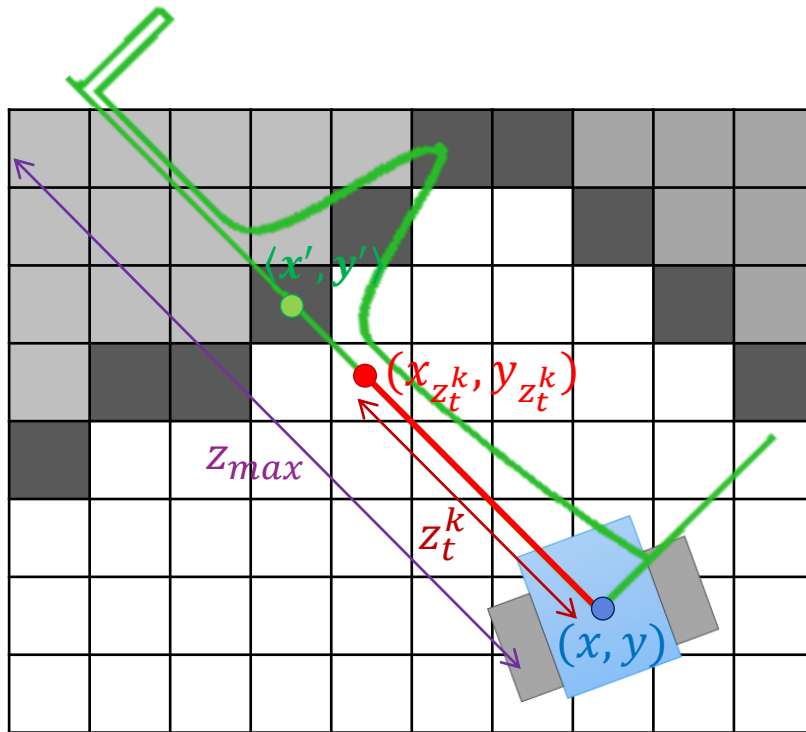
(a) Sonar data



(b) Laser data



# Likelihood Field



- 1: Algorithm `likelihood_field_range_finder_model( $z_t, x_t, m$ )`:
- 2:      $q = 1$
- 3:     for all  $k$  do
- 4:         if  $z_t^k \neq z_{max}$
- 5:              $x_{z_t^k} = x + x_{k,sens} \cos \theta - y_{k,sens} \sin \theta + z_t^k \cos(\theta + \theta_{k,sens})$
- 6:              $y_{z_t^k} = y + y_{k,sens} \cos \theta + x_{k,sens} \sin \theta + z_t^k \sin(\theta + \theta_{k,sens})$
- 7:              $dist = \min_{x', y'} \left\{ \sqrt{(x_{z_t^k} - x')^2 + (y_{z_t^k} - y')^2} \mid \langle x', y' \rangle \text{ occupied in } m \right\}$
- 8:              $q = q \cdot (z_{hit} \cdot \text{prob}(dist, \sigma_{hit}) + \frac{z_{random}}{z_{max}})$
- 9:     return  $q$

# Grid-based Fast-SLAM

1. Predict the next pose  $x_t^{(i)}$  by motion model.

$$x_t^{(i)} \sim p(x_t^{(i)} | x_{t-1}^{(i)}, u_{t-1})$$

2. Update the occupancy grid map of each particle.

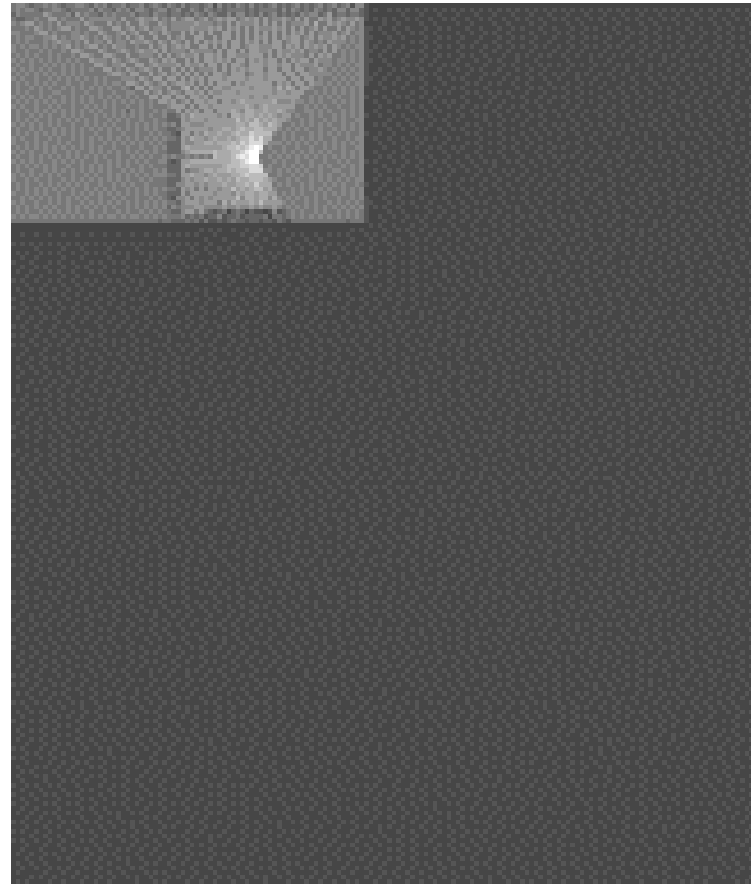
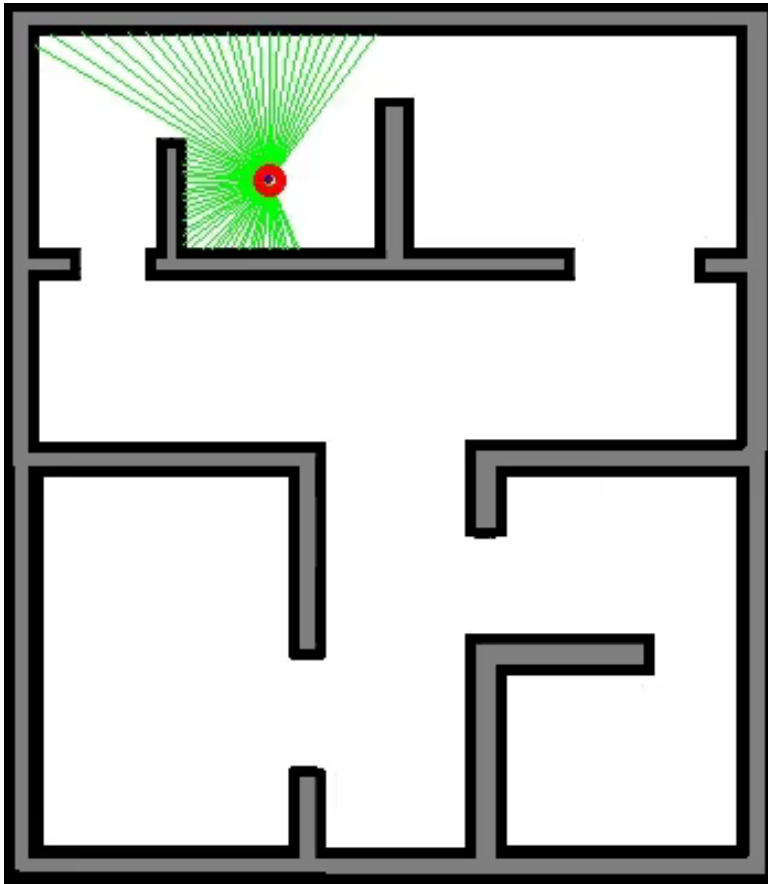
$$\log Odd(s|z) = \log \frac{p(z|s=1)}{p(z|s=0)} + \log Odd(s)$$

3. Update the importance weight of particles.

$$w_t^{(i)} = \eta \prod_i (z_{hit} \cdot prob(dist, \sigma_{hit}) + \frac{z_{random}}{z_{max}})$$

4. Resampling.

## Grid-based Fast-SLAM Demo



# Outline

- State Estimation and SLAM Problem
- SLAM Back-end (Error Compensation)
  - Filter-based Methods
    - Probability Theory and Bayes Filter
    - Kalman Filter (KF) / Extended Kalman Filter (EKF)
      - EKF-SLAM
    - Particle Filter
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  - Graph-based Methods
    - Pose Graph and Least-square Optimization
    - Gauss-Newton and Levenberg-Marquardt Algorithm
    - Sparse Matrix for Optimization



# State Estimation

$$\begin{aligned} \mathbf{x} \sim \mathcal{N}(\boldsymbol{\mu}, \boldsymbol{\Sigma}) \quad P(\mathbf{x}) &= \frac{1}{\sqrt{(2\pi)^N \det(\boldsymbol{\Sigma})}} \exp\left(-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu})\right) \\ -\ln P(\mathbf{x}) &= \frac{1}{2} \ln((2\pi)^N \det(\boldsymbol{\Sigma})) + \frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu}) \end{aligned}$$

$$\mathbf{x}_k = f(\mathbf{x}_{k-1}, \mathbf{u}_k) + \mathbf{w}_k, \quad \mathbf{w}_k \sim \mathcal{N}(0, \mathbf{R}_k)$$

$$\mathbf{z}_{k,j} = h(\mathbf{m}_j, \mathbf{x}_k) + \mathbf{v}_{k,j}, \quad \mathbf{v}_{k,j} \sim \mathcal{N}(0, \mathbf{Q}_{k,j})$$

- Probability of  $\{\mathbf{x}_1, \dots, \mathbf{x}_N, \mathbf{m}_1, \dots, \mathbf{m}_M\}$  given  $\{\mathbf{u}_1, \dots, \mathbf{u}_N, \mathbf{z}_{1,1}, \dots, \mathbf{z}_{N,M}\}$  :

$$P(\mathbf{x}, \mathbf{m} | \mathbf{z}, \mathbf{u}) = \frac{P(\mathbf{z}, \mathbf{u} | \mathbf{x}, \mathbf{m}) P(\mathbf{x}, \mathbf{m})}{P(\mathbf{z}, \mathbf{u})} \propto \underbrace{P(\mathbf{z}, \mathbf{u} | \mathbf{x}, \mathbf{m})}_{\text{likelihood}} \underbrace{P(\mathbf{x}, \mathbf{m})}_{\text{prior}}$$

$$(\mathbf{x}, \mathbf{m})_{MAP}^* = \operatorname{argmax} P(\mathbf{x}, \mathbf{m} | \mathbf{z}, \mathbf{u}) = \operatorname{argmax} P(\mathbf{z}, \mathbf{u} | \mathbf{x}, \mathbf{m}) P(\mathbf{x}, \mathbf{m})$$

$$(\mathbf{x}, \mathbf{m})_{MLE}^* = \operatorname{argmax} P(\mathbf{z}, \mathbf{u} | \mathbf{x}, \mathbf{m})$$

$$P(\mathbf{z}_{k,j} | \mathbf{x}_k, \mathbf{m}_j) = \mathcal{N}(h(\mathbf{m}_j, \mathbf{x}_k), \mathbf{Q}_{k,j})$$

$$(\mathbf{x}_k, \mathbf{m}_j)_{MLE}^* = \operatorname{argmax} \mathcal{N}(h(\mathbf{m}_j, \mathbf{x}_k), \mathbf{Q}_{k,j}) = \operatorname{argmin} \frac{1}{2} (\mathbf{z}_{k,j} - h(\mathbf{m}_j, \mathbf{x}_k))^T \mathbf{Q}_{k,j}^{-1} (\mathbf{z}_{k,j} - h(\mathbf{m}_j, \mathbf{x}_k))$$

# State Estimation

$$(\mathbf{x}_k, \mathbf{m}_j)_{MLE}^* = \operatorname{argmin} \frac{1}{2} \left( \mathbf{z}_{k,j} - h(\mathbf{m}_j, \mathbf{x}_k) \right)^T \mathbf{Q}_{k,j}^{-1} \left( \mathbf{z}_{k,j} - h(\mathbf{m}_j, \mathbf{x}_k) \right)$$

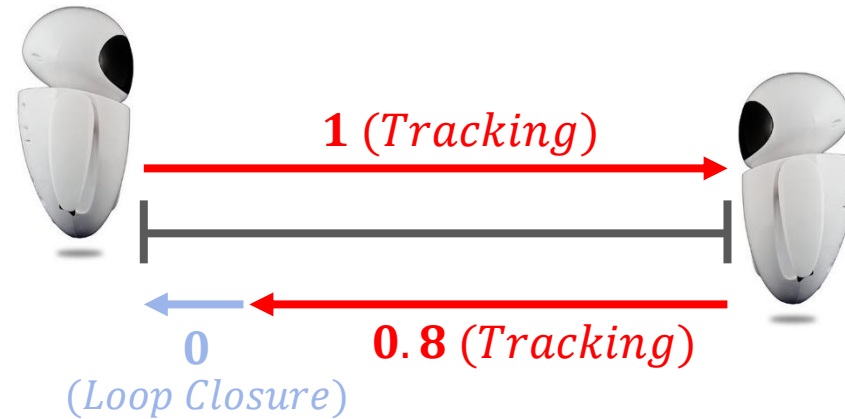
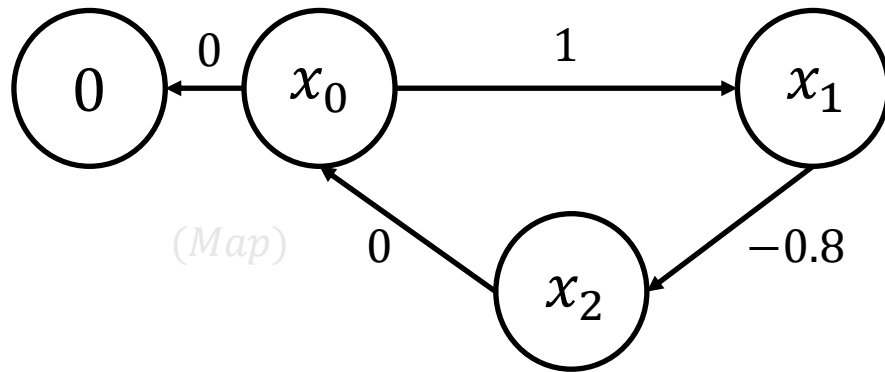
$$P(\mathbf{z}, \mathbf{u} | \mathbf{x}, \mathbf{m}) = \prod_k P(\mathbf{u}_k | \mathbf{x}_{k-1}, \mathbf{x}_k) \prod_{k,j} P(\mathbf{z}_{k,j} | \mathbf{x}_k, \mathbf{m}_j)$$

$$\mathbf{e}_{\mathbf{u},k} = \mathbf{x}_k - f(\mathbf{x}_{k-1}, \mathbf{u}_k)$$

$$\mathbf{e}_{\mathbf{z},k,j} = \mathbf{z}_{k,j} - h(\mathbf{m}_j, \mathbf{x}_k)$$

$$\min F(\mathbf{x}, \mathbf{m}) = \sum_k \mathbf{e}_{\mathbf{u},k}^T \mathbf{R}_K^{-1} \mathbf{e}_{\mathbf{u},k} + \sum_k \sum_j \mathbf{e}_{\mathbf{z},k,j}^T \mathbf{Q}_{K,j}^{-1} \mathbf{e}_{\mathbf{z},k,j}$$

# Graph Optimization: 1D Example



Error function

$$x_0 = 0$$

$$x_1 = x_0 + 1$$

$$x_2 = x_1 - 0.8$$

$$x_0 = x_2 + 0$$



$$f_1 = x_0$$

$$f_2 = x_1 - x_0 - 1$$

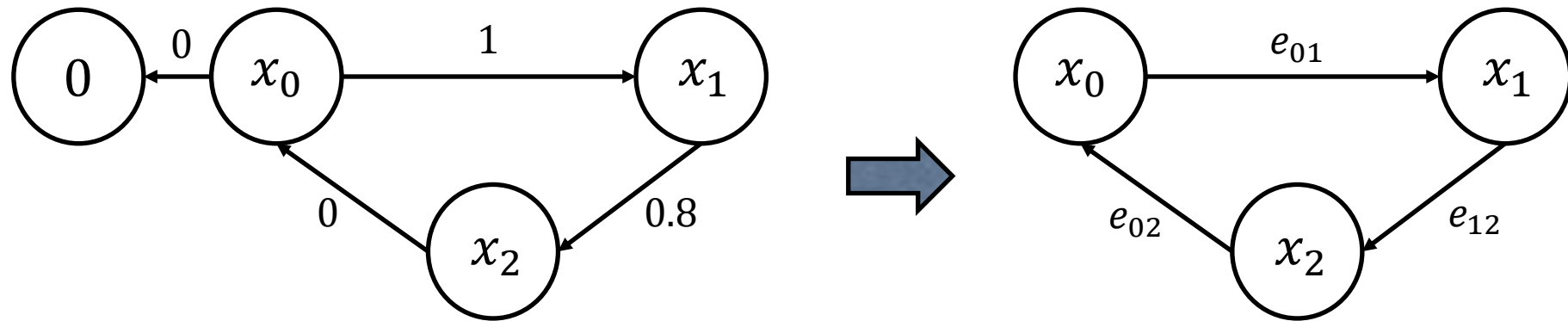
$$f_3 = x_2 - x_1 + 0.8$$

$$f_4 = x_0 - x_2$$

$$\min_x \sum_i w_i f_i^2 = w_1 x_0^2 + w_2 (x_1 - x_0 - 1)^2 + w_3 (x_2 - x_1 + 0.8)^2 + w_4 (x_0 - x_2)^2$$

(Optimization)

## Graph Optimization: 1D Example



### Error Function

$$e_{01} = x_1 - x_0 - 1$$

$$e_{12} = x_2 - x_1 - 0.8$$

$$e_{02} = x_0 - x_2$$

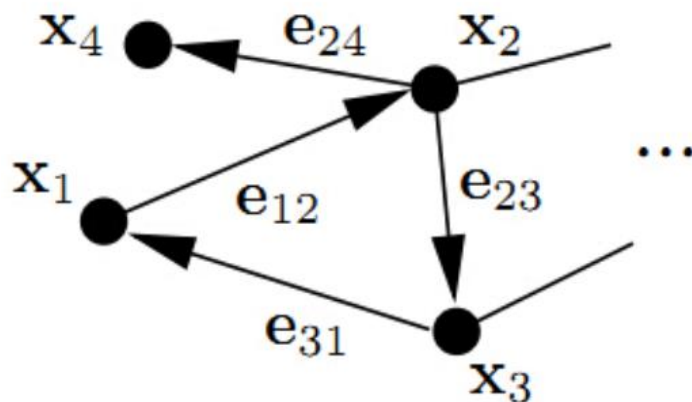
$$\min_x \sum_{i,j} w_{ij} e_{ij}^2 = w_{01}(x_1 - x_0 - 1)^2 + w_{12}(x_2 - x_1 + 0.8)^2 + w_{02}(x_0 - x_2)^2$$

## Graph Optimization: General Form

$$\min_x \sum_{i,j} w_{ij} e_{ij}^2 = w_{01}(x_1 - x_0 - 1)^2 + w_{12}(x_2 - x_1 + 0.8)^2 + w_{02}(x_0 - x_2)^2$$

$$\mathbf{F}(\mathbf{x}) = \sum_{\langle i,j \rangle \in \mathcal{C}} \underbrace{\mathbf{e}(\mathbf{x}_i, \mathbf{x}_j, \mathbf{z}_{ij})^\top \boldsymbol{\Omega}_{ij} \mathbf{e}(\mathbf{x}_i, \mathbf{x}_j, \mathbf{z}_{ij})}_{\mathbf{F}_{ij}} \quad (1)$$



$$\mathbf{x}^* = \underset{\mathbf{x}}{\operatorname{argmin}} \mathbf{F}(\mathbf{x}). \quad (2)$$



$$\begin{aligned} \mathbf{F}(\mathbf{x}) = & \mathbf{e}_{12}^\top \boldsymbol{\Omega}_{12} \mathbf{e}_{12} \\ & + \mathbf{e}_{23}^\top \boldsymbol{\Omega}_{23} \mathbf{e}_{23} \\ & + \mathbf{e}_{31}^\top \boldsymbol{\Omega}_{31} \mathbf{e}_{31} \\ & + \mathbf{e}_{24}^\top \boldsymbol{\Omega}_{24} \mathbf{e}_{24} \\ & + \dots \end{aligned}$$

# Graph Optimization Library

## g2o - General Graph Optimization

Linux:   Windows:  

g2o is an open-source C++ framework for optimizing graph-based nonlinear error functions. g2o has been designed to be easily extensible to a wide range of problems and a new problem typically can be specified in a few lines of code. The current implementation provides solutions to several variants of SLAM and BA.

<https://github.com/RainerKuemmerle/g2o>

## Ceres Solver

Ceres Solver is an open source C++ library for modeling and solving large, complicated optimization problems. It is a feature rich, mature and performant library which has been used in production at Google since 2010. Ceres Solver can solve two kinds of problems.

<https://github.com/ceres-solver/ceres-solver>

# Non-linear Optimization

# Basics of Optimization

## Least Squares Problem

Find  $\mathbf{x}^*$ , a local minimizer for

$$F(\mathbf{x}) = \frac{1}{2} \sum_{i=1}^m (f_i(\mathbf{x}))^2 ,$$

where  $f_i : \mathbb{R}^n \mapsto \mathbb{R}$ ,  $i = 1, \dots, m$  are given functions, and  $m \geq n$ .

$m$ : number of data points

$n$ : number of parameters

$$\frac{dF}{d\mathbf{x}} = 0$$

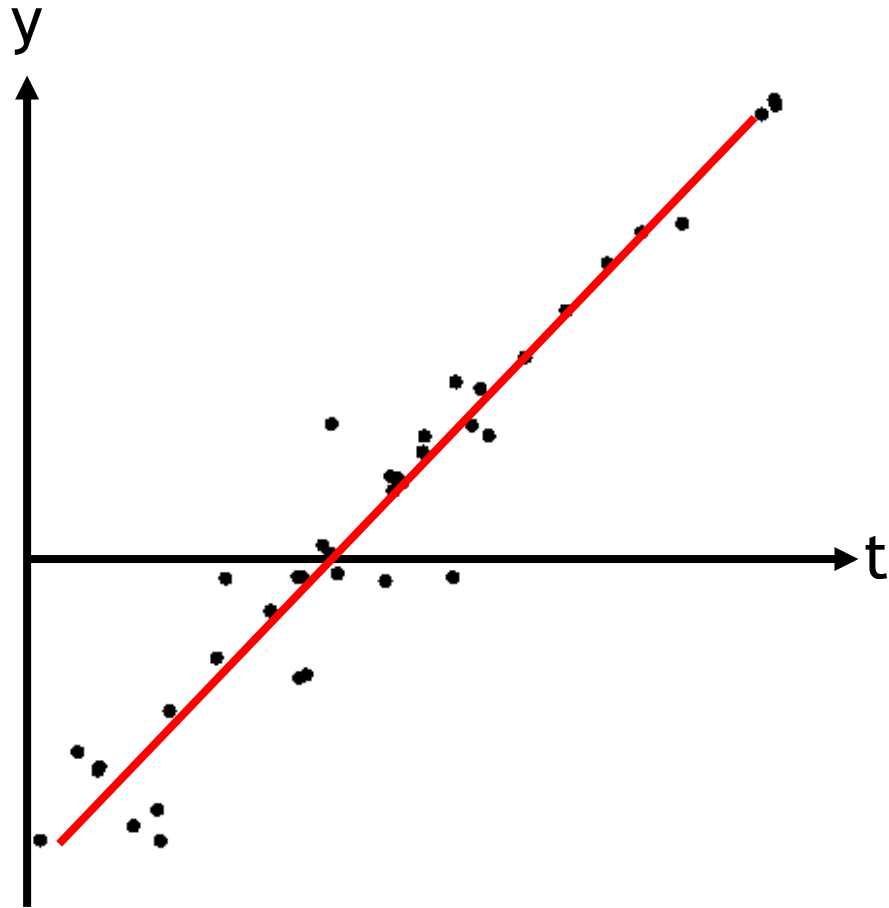
## Local Minimizer

Given  $F : \mathbb{R}^n \mapsto \mathbb{R}$ . Find  $\mathbf{x}^*$  so that

$$F(\mathbf{x}^*) \leq F(\mathbf{x}) \quad \text{for} \quad \|\mathbf{x} - \mathbf{x}^*\| < \delta .$$



## Example: Linear Least Square Fitting



model

parameters

$$y(t) = \boxed{M}(t; \boxed{\mathbf{x}}) = x_0 + x_1 t$$

$$f_i(x) = y_i - \boxed{M(t_i; \mathbf{x})}$$

Residual(error)

prediction

$M(t; \mathbf{x}) = x_0 + x_1 t + x_2 t^3$  is linear, too.

# Example: Nonlinear Least Square Fitting

parameters

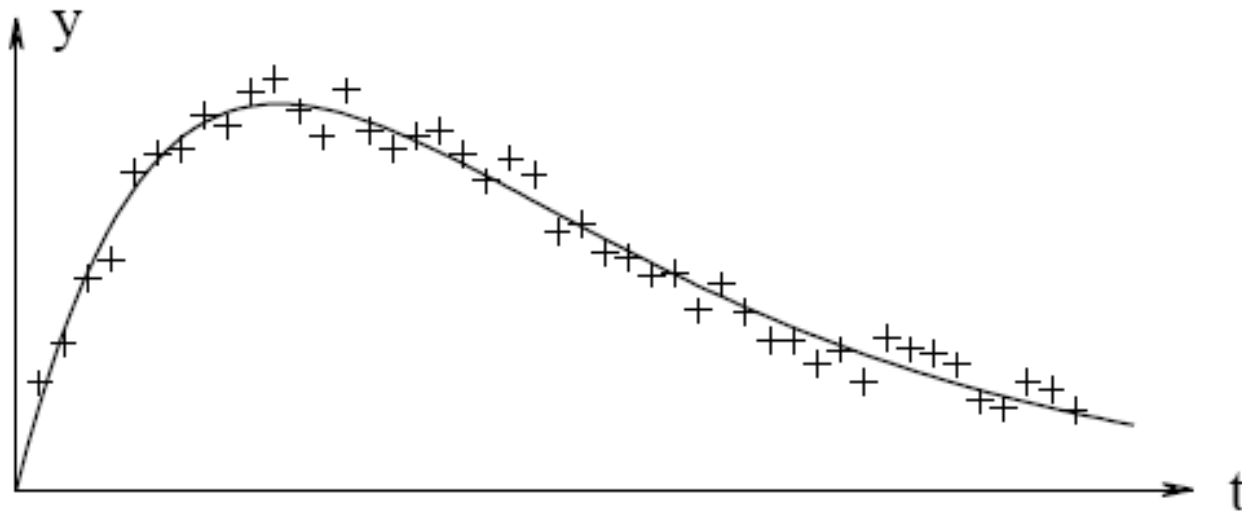
$$\mathbf{x} = [x_1, x_2, x_3, x_4]^T$$

model

$$M(t; \mathbf{x}) = x_3 e^{x_1 t} + x_4 e^{x_2 t}$$

residuals

$$\begin{aligned} f_i(\mathbf{x}) &= y_i - M(t_i; \mathbf{x}) \\ &= y_i - (x_3 e^{x_1 t} + x_4 e^{x_2 t}) \end{aligned}$$



# Function Minimization

*Taylor expansion*  $F(\mathbf{x} + \mathbf{h}) \approx F(\mathbf{x}) + \mathbf{J}(\mathbf{x})^T \mathbf{h} + \frac{1}{2} \mathbf{h}^T \mathbf{H}(\mathbf{x}) \mathbf{h}$

$$\mathbf{J}(\mathbf{x}) \equiv \mathbf{F}'(\mathbf{x}) = \begin{bmatrix} \frac{\partial F}{\partial x_1}(\mathbf{x}) \\ \vdots \\ \frac{\partial F}{\partial x_n}(\mathbf{x}) \end{bmatrix}$$

$$\mathbf{H}(\mathbf{x}) \equiv \mathbf{F}''(\mathbf{x}) = \begin{bmatrix} \frac{\partial^2 F}{\partial x_1^2}(\mathbf{x}) & \frac{\partial^2 F}{\partial x_1 \partial x_2}(\mathbf{x}) & \dots & \frac{\partial^2 F}{\partial x_1 \partial x_n}(\mathbf{x}) \\ \frac{\partial^2 F}{\partial x_2 \partial x_1}(\mathbf{x}) & \frac{\partial^2 F}{\partial x_2^2}(\mathbf{x}) & \dots & \frac{\partial^2 F}{\partial x_2 \partial x_n}(\mathbf{x}) \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial^2 F}{\partial x_n \partial x_1}(\mathbf{x}) & \frac{\partial^2 F}{\partial x_n \partial x_2}(\mathbf{x}) & \dots & \frac{\partial^2 F}{\partial x_n^2}(\mathbf{x}) \end{bmatrix}$$

# Function Minimization

Necessary condition for a local minimizer :

$$J(\mathbf{x}^*) \equiv F'(\mathbf{x}) = \mathbf{0}$$

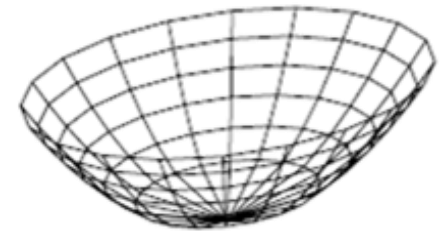
Why?

By definition, if  $\mathbf{x}^*$  is a local minimizer,

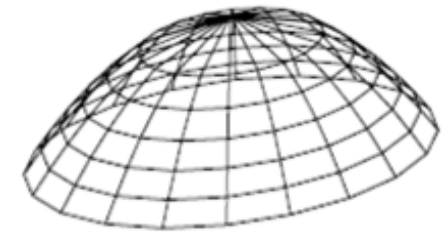
$$\|\mathbf{h}\| \text{ is small enough } \rightarrow F(\|\mathbf{x}^* + \mathbf{h}\|) > F(\mathbf{x}^*)$$

$$F(\mathbf{x}^* + \mathbf{h}) \approx F(\mathbf{x}^*) + J(\mathbf{x}^*)^T \mathbf{h} + \frac{1}{2} \mathbf{h}^T H(\mathbf{x}) \mathbf{h} > F(\mathbf{x}^*)$$

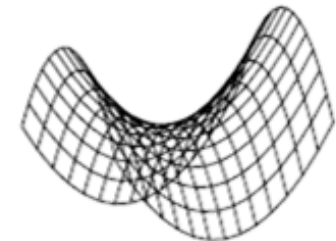
$$F(\mathbf{x}^* - \mathbf{h}) \approx F(\mathbf{x}^*) - J(\mathbf{x}^*)^T \mathbf{h} + \frac{1}{2} \mathbf{h}^T H(\mathbf{x}) \mathbf{h} < F(\mathbf{x}^*)$$



a) *minimum*



b) *maximum*



c) *saddle point*

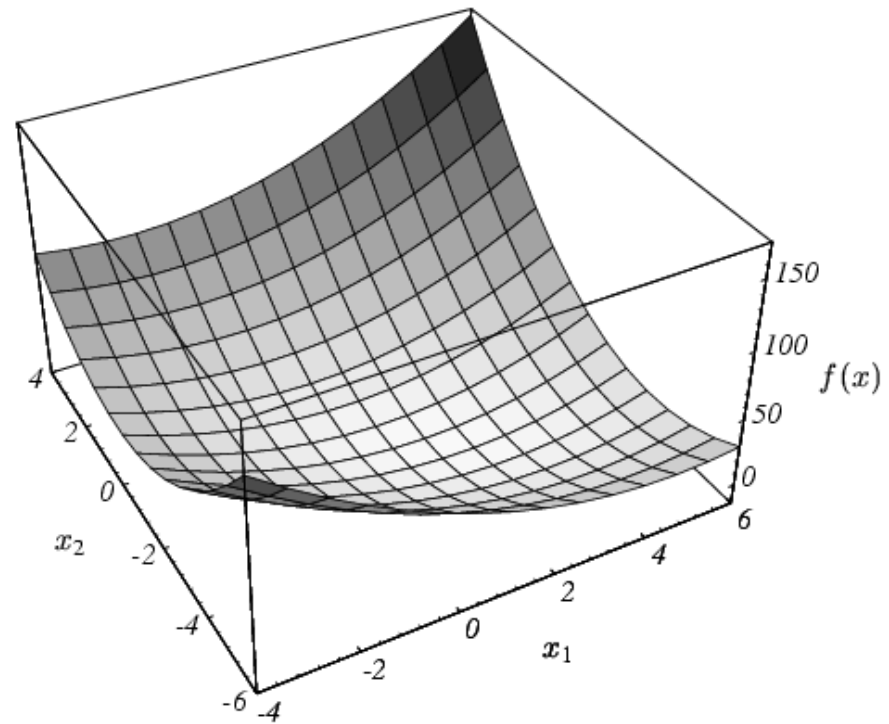
# Quadratic Functions

$$F(\mathbf{x}) = \frac{1}{2} \mathbf{x}^T \mathbf{A} \mathbf{x} - \mathbf{b}^T \mathbf{x} + \mathbf{c}$$

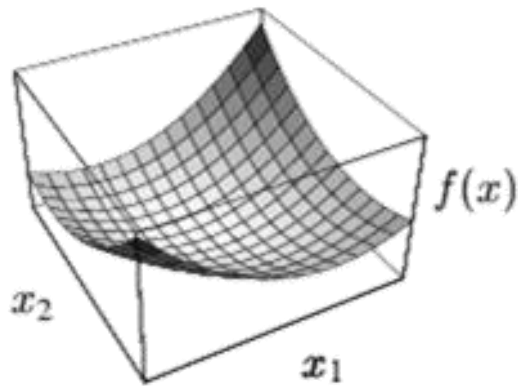
$$\mathbf{A} = \begin{bmatrix} 3 & 2 \\ 2 & 6 \end{bmatrix}$$

$$\mathbf{b} = \begin{bmatrix} 2 \\ -8 \end{bmatrix}$$

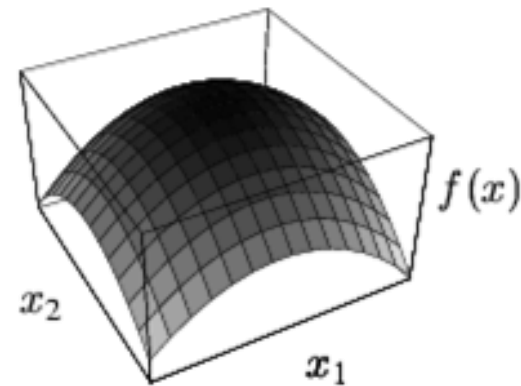
$$\mathbf{c} = 0$$



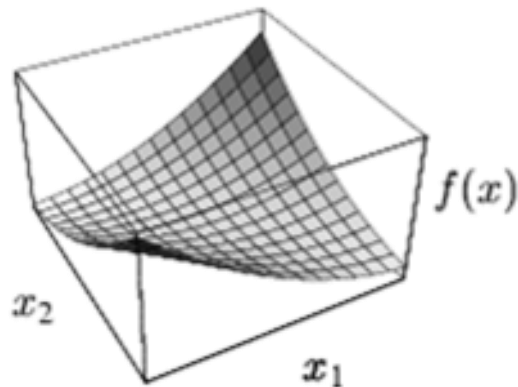
# Quadratic Functions



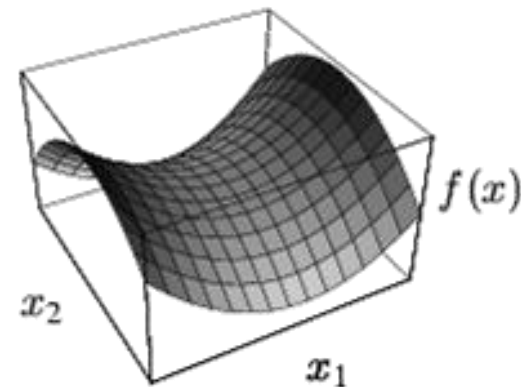
$\mathbf{A}$  is **positive definite**.  
All eigenvalues are positive.  
For all  $x$ ,  $x^T \mathbf{A} x > 0$ .



$\mathbf{A}$  is **negative definite**.  
All eigenvalues are negative.  
For all  $x$ ,  $x^T \mathbf{A} x < 0$ .



$\mathbf{A}$  is **singular**



$\mathbf{A}$  is **indefinite**

# Descent Methods

$\mathbf{x}_0, \mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_k \rightarrow \mathbf{x}^*$  for  $k \rightarrow \infty$

1. Find a descent direction  $\mathbf{h}_d$
2. find a step length giving a good decrease in the  $F$ -value.

## Algorithm Descent method

**begin**

$k := 0; \mathbf{x} := \mathbf{x}_0; found := \mathbf{false}$  {Starting point}

**while** (**not** *found*) **and** ( $k < k_{\max}$ )

$\mathbf{h}_d := \text{search\_direction}(\mathbf{x})$  {From  $\mathbf{x}$  and downhill}

**if** (no such  $\mathbf{h}$  exists)

$found := \mathbf{true}$  { $\mathbf{x}$  is stationary}

**else**

$\alpha := \text{step\_length}(\mathbf{x}, \mathbf{h}_d)$  {from  $\mathbf{x}$  in direction  $\mathbf{h}_d$ }

$\mathbf{x} := \mathbf{x} + \alpha \mathbf{h}_d; k := k+1$  {next iterate}

**end**

## Descent Direction (Line Search Method)

$$\begin{aligned} F(\mathbf{x} + \alpha \mathbf{h}) &= F(\mathbf{x}) + \alpha \mathbf{h}^\top \mathbf{F}'(\mathbf{x}) + O(\alpha^2) \\ &\simeq F(\mathbf{x}) + \alpha \mathbf{h}^\top \mathbf{F}'(\mathbf{x}) \quad \text{for } \alpha \text{ sufficiently small.} \end{aligned}$$

Definition of descent direction:

$\mathbf{h}$  is a descent direction for  $F$  at  $\mathbf{x}$  if  $\mathbf{h}^\top \mathbf{F}'(\mathbf{x}) < 0$



## Steepest Descent Method

$$\begin{aligned} F(\mathbf{x} + \alpha \mathbf{h}) &= F(\mathbf{x}) + \alpha \mathbf{h}^\top \mathbf{F}'(\mathbf{x}) + O(\alpha^2) \\ &\simeq F(\mathbf{x}) + \alpha \mathbf{h}^\top \mathbf{F}'(\mathbf{x}) \quad \text{for } \alpha \text{ sufficiently small.} \end{aligned}$$

$$\boxed{\frac{F(\mathbf{x}) - F(\mathbf{x} + \alpha \mathbf{h})}{\alpha \|\mathbf{h}\|}} = -\frac{1}{\|\mathbf{h}\|} \mathbf{h}^\top \mathbf{F}'(\mathbf{x}) = -\|\mathbf{F}'(\mathbf{x})\| \cos \theta$$

the decrease of  $F(\mathbf{x})$  per unit along  $\mathbf{h}$  direction

greatest gain rate if  $\theta = \pi \rightarrow \mathbf{h}_{\text{sd}} = -\mathbf{F}'(\mathbf{x})$

$\mathbf{h}_{\text{sd}}$  is a descent direction because  $\mathbf{h}_{\text{sd}}^\top \mathbf{F}'(\mathbf{x}) = -F'(\mathbf{x})^2 < 0$

# Steepest Descent Method

$\varphi(\alpha) = F(\mathbf{x} + \alpha \mathbf{h})$ ,  $\mathbf{x}$  and  $\mathbf{h}$  are fixed,  $\alpha \geq 0$ .

Find  $\alpha$  so that  $\varphi(\alpha) = F(\mathbf{x} + \alpha \mathbf{h})$  is minimum.

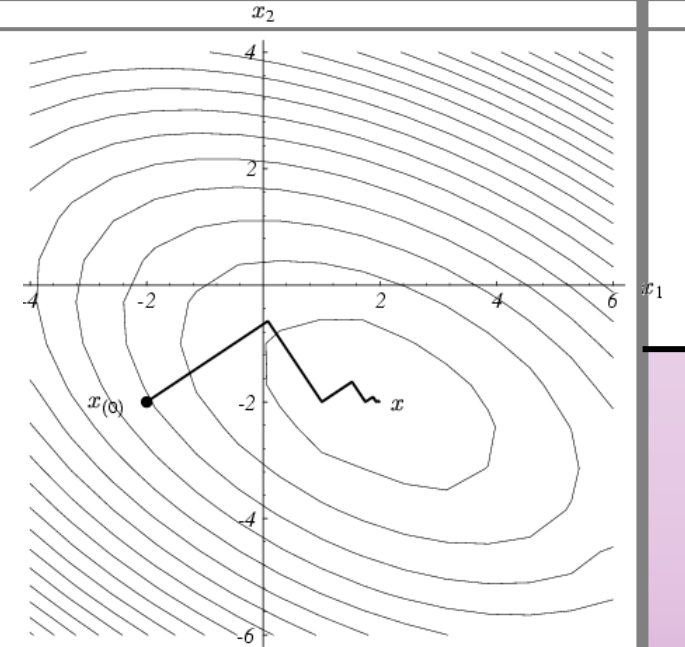
$$0 = \frac{\partial \varphi(\alpha)}{\partial \alpha} = \frac{\partial F(\mathbf{x} + \alpha \mathbf{h})}{\partial \alpha} = \frac{\partial F(\mathbf{x} + \alpha \mathbf{h})}{\partial (\mathbf{x} + \alpha \mathbf{h})} \frac{\partial (\mathbf{x} + \alpha \mathbf{h})}{\partial \alpha} = \mathbf{h}^T F'(\mathbf{x} + \alpha \mathbf{h})$$

$$\mathbf{h} = -F'(\mathbf{x})$$

$$= \mathbf{h}^T (F'(\mathbf{x}) + \alpha F''(\mathbf{x})^T \mathbf{h}) = \mathbf{h}^T (-\mathbf{h} + \alpha \mathbf{H} \mathbf{h})$$

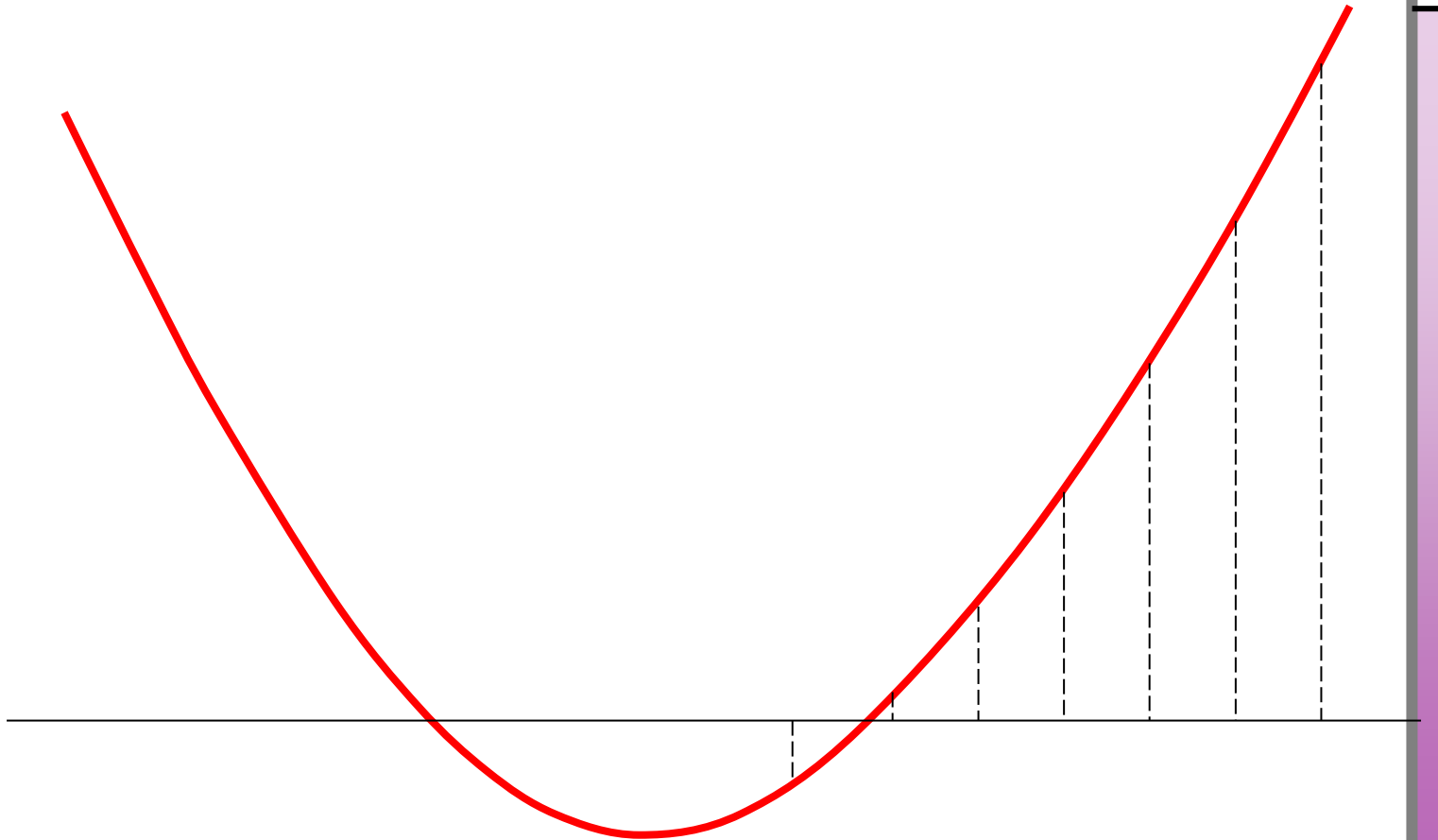
$$\alpha = \frac{\mathbf{h}^T \mathbf{h}}{\mathbf{h}^T \mathbf{H} \mathbf{h}}$$

Problem: Has good performance in the initial stages of the iterative process, but converge very slow with a linear rate.



# Newton's Method

- Root finding for  $f(x)=0$
- March  $x$  and test signs
- Determine  $\Delta x$   
(small  $\rightarrow$  slow; large  $\rightarrow$  miss)



# Newton's Method

- Root finding for  $f(x)=0$

**Taylor's expansion:**

$$f(x_0 + \varepsilon) = f(x_0) + f'(x_0)\varepsilon + \frac{1}{2}f''(x_0)\varepsilon^2 + \dots$$

$$0 = f(x_0 + \varepsilon) \approx f(x_0) + f'(x_0)\varepsilon$$

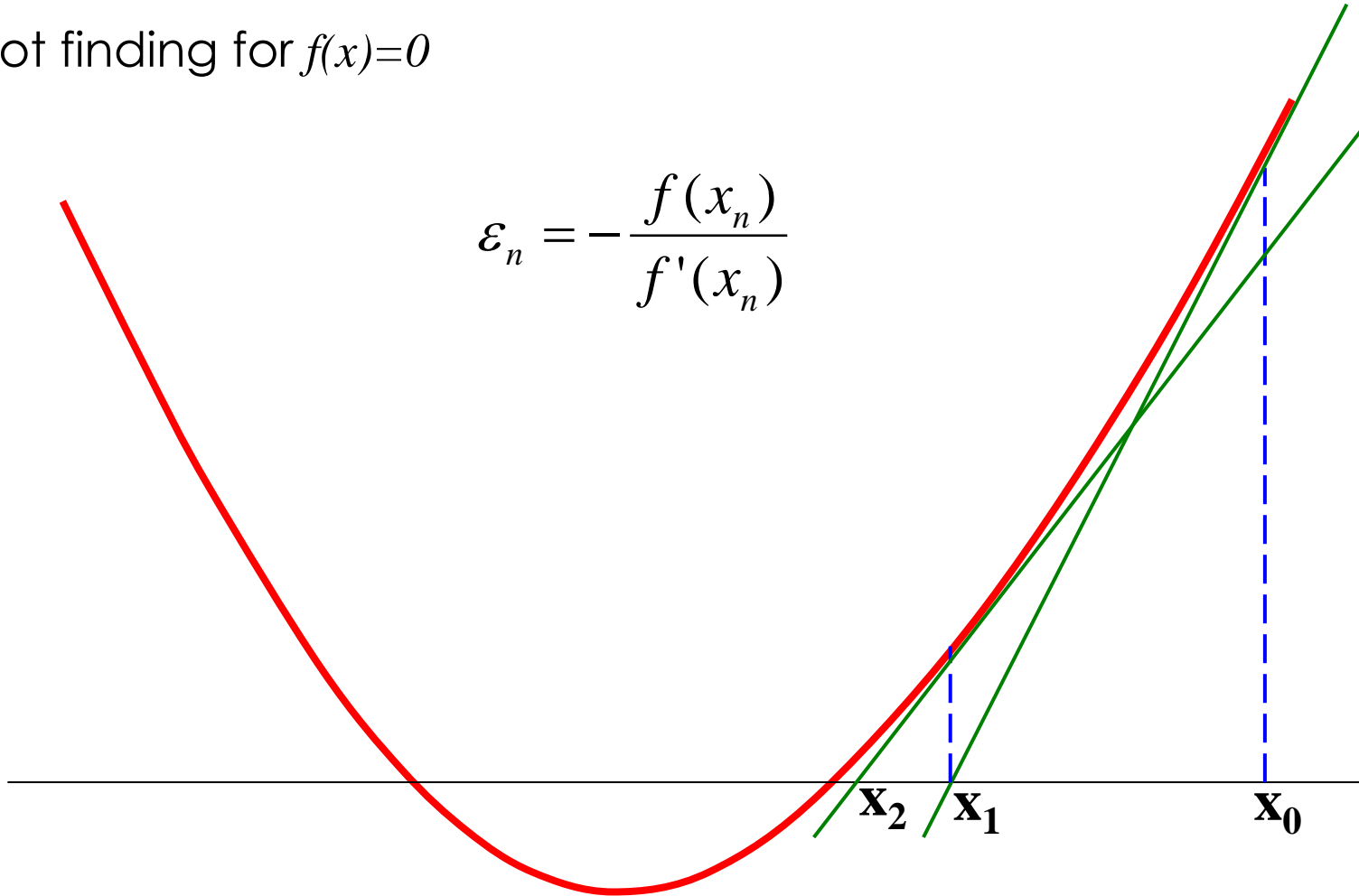
$$\varepsilon = -\frac{f(x_0)}{f'(x_0)}$$

$$x_{n+1} = x_n - \frac{f(x_n)}{f'(x_n)}$$

# Newton's Method

- Root finding for  $f(x)=0$

$$\varepsilon_n = -\frac{f(x_n)}{f'(x_n)}$$



## Newton's Method

$\mathbf{x}^*$  is a stationary point  $\rightarrow$  it satisfies  $\mathbf{F}'(\mathbf{x}^*) = \mathbf{0}$ .

$$\begin{aligned}\mathbf{F}'(\mathbf{x}+\mathbf{h}) &= \mathbf{F}'(\mathbf{x}) + \mathbf{F}''(\mathbf{x})\mathbf{h} + O(\|\mathbf{h}\|^2) \\ &\simeq \mathbf{F}'(\mathbf{x}) + \mathbf{F}''(\mathbf{x})\mathbf{h} \quad \text{for } \|\mathbf{h}\| \text{ sufficiently small} \\ &= \mathbf{0}\end{aligned}$$

$$\rightarrow \mathbf{H} \mathbf{h}_n = -\mathbf{F}'(\mathbf{x}) \quad \text{with } \mathbf{H} = \mathbf{F}''(\mathbf{x})$$

$$\mathbf{x} := \mathbf{x} + \mathbf{h}_n$$

Suppose that  $\mathbf{H}$  is positive definite

$$\rightarrow \mathbf{u}^\top \mathbf{H} \mathbf{u} > 0 \text{ for all nonzero } \mathbf{u}.$$

$$\rightarrow 0 < \mathbf{h}_n^\top \mathbf{H} \mathbf{h}_n = -\mathbf{h}_n^\top \mathbf{F}'(\mathbf{x})$$

$$\rightarrow \mathbf{h}_n \text{ is a descent direction}$$

# Newton's Method

$$\mathbf{H}\mathbf{h} = -F'(\mathbf{x})$$

$$\mathbf{h} = -\mathbf{H}^{-1}\mathbf{J}$$

- It has good performance in the final stage of the iterative process, where  $\mathbf{x}$  is close to  $\mathbf{x}^*$ .
- It requires solving a linear system and  $\mathbf{H}$  is not always positive definite.

→ Use the approximate Hessian  $\mathbf{H} \approx \mathbf{J}^T \mathbf{J}$  Gauss-Newton

# Gauss-Newton

$$\mathbf{h}^* = \operatorname{argmin} \frac{1}{2} \sum_{i=1}^m \|f_i(\mathbf{x} + \mathbf{h})\|^2$$

$$f(\mathbf{x} + \mathbf{h}) \approx f(\mathbf{x}) + \mathbf{J}(\mathbf{x})^T \mathbf{h}$$

$$\frac{1}{2} \|f(\mathbf{x} + \mathbf{h})\|^2 \approx \frac{1}{2} \|f(\mathbf{x}) + \mathbf{J}(\mathbf{x})^T \mathbf{h}\|^2 = \frac{1}{2} (\|f(\mathbf{x})\|^2 + 2f(\mathbf{x})\mathbf{J}(\mathbf{x})^T \mathbf{h} + \mathbf{h}^T \mathbf{J}(\mathbf{x})\mathbf{J}(\mathbf{x})^T \mathbf{h})$$

$$\mathbf{J}(\mathbf{x})f(\mathbf{x})^T + \mathbf{J}(\mathbf{x})\mathbf{J}(\mathbf{x})^T \mathbf{h} = \mathbf{0}$$

$$\underbrace{\mathbf{J}(\mathbf{x})\mathbf{J}(\mathbf{x})^T}_{\mathbf{H}(\mathbf{x})} \mathbf{h} = -\underbrace{\mathbf{J}(\mathbf{x})f(\mathbf{x})^T}_{\mathbf{g}(\mathbf{x})}$$

Newton's Method:

$$\mathbf{H}\mathbf{h} = -F'(\mathbf{x})$$



# Levenberg-Marquardt Method (LM)

- LM can be thought of as a combination of steepest descent and the Newton method.
  - When the current solution is far from the correct one, the algorithm behaves like a steepest descent method: slow, but guaranteed to converge.
  - When the current solution is close to the correct solution, it becomes a Newton's method.

**if**  $F''(\mathbf{x})$  is positive definite  
     $\mathbf{h} := \mathbf{h}_n$   
**else**  
     $\mathbf{h} := \mathbf{h}_{sd}$   
     $\mathbf{x} := \mathbf{x} + \alpha \mathbf{h}$

true-region method

$$\rho = \frac{f(\mathbf{x} + \mathbf{h}) - f(\mathbf{x})}{\mathbf{J}(\mathbf{x})^T \mathbf{h}}$$

This needs to calculate second-order derivative which might not be available.

# Levenberg-Marquardt Method (LM)

Initialize  $\mathbf{x} = \mathbf{x}_0, \mu = \mu_0$

For  $i=0 \sim K$

Find  $\mathbf{h}$  such that  $\min_{\mathbf{h}} \frac{1}{2} \|\mathbf{f}(\mathbf{x}) + \mathbf{J}(\mathbf{x})^T \mathbf{h}\|^2$  s.t.  $\|\mathbf{D}\mathbf{h}\|^2 \leq \mu$

Calculate  $\rho$

If  $\rho \geq \frac{3}{4}$

$$\mu = 2\mu$$

If  $\rho < \frac{1}{4}$

$$\mu = 0.5\mu$$

If  $\rho \geq Th$

else  $\mathbf{x} = \mathbf{x} + \mathbf{h}$

If  $\mathbf{h}$  is smaller than  $\epsilon$ , stop

$$\mathcal{L}(\mathbf{h}, \lambda) = \frac{1}{2} \|\mathbf{f}(\mathbf{x}) + \mathbf{J}(\mathbf{x})^T \mathbf{h}\|^2 + \lambda (\|\mathbf{D}\mathbf{h}\|^2 - \mu)$$

$$\nabla \mathcal{L}(\mathbf{h}, \lambda) = 0$$

$$(\mathbf{J}(\mathbf{x})\mathbf{J}(\mathbf{x})^T + \lambda \mathbf{D}^T \mathbf{D}) \mathbf{h} = -\mathbf{J}(\mathbf{x}) \mathbf{f}(\mathbf{x})^T$$

$$(\mathbf{H}(\mathbf{x}) + \lambda \mathbf{I}) \mathbf{h} = -\mathbf{g}(\mathbf{x})$$