


# Grid Maps

1. What it is

Grid map is a way to represent the map. World is divided into grid of cells   
We can use this to draw a map (large map may need large memory).

- (2) Assumption 1: each cell is totally occupied or totally free
- (3) Assumption 2: world is static. Cells don't change state. (no true in reality)
- (4) Assumption 3: cells are independent (not very true in reality).

## 2. Representation.

naturally we use a binary random variable to model a cell:

$p(m_i)=1$  occupied.

$p(m_i)=0$  not occupied.

but we also allow  $p(m_i)=0.5$  not known

Given sensor data.

e.g. 

$m_1$	$m_2$
$m_3$	$m_4$

  
 $m_1=0.9$   
 $m_2=0.5$   
 $m_3=0.8$   
 $m_4=0.1$

Probability of a state.

$$P(M = \begin{bmatrix} 0.9 \\ 0.5 \\ 0.8 \\ 0.1 \end{bmatrix}) = \prod_{i=1}^4 p(m_i = m_i) = 0.9 \times (1-0.5) \times 0.8 \times (1-0.1)$$

## 3. Mapping with known poses.

The problem: Given sensor data and pose of sensor (robot), estimate a map.

Mathematically:

$$P(m | z_{1:t}, x_{1:t}) = \prod_i p(m_i | z_{1:t}, x_{1:t})$$

Idea 1: To deal with binary variable, we can compute ratio  $\frac{p(z_t)}{1-p(z_t)}$  for estimation.

Idea 2: Using Bayes Recursive Filter. That is, find relation between current state and previous state, thus form a recursive algorithm

Go for  $p(m_i | z_{1:t+1}, x_{1:t+1})$ .  
knowing  $x_t$ , past pose of robot does not help

$$p(m_i | z_{1:t}, x_{1:t}) \xrightarrow{\text{Bayes Rule}} \frac{p(z_t | m_i, z_{1:t-1}, x_{1:t}) p(m_i | z_{1:t-1}, x_{1:t-1})}{p(z_t | z_{1:t-1}, x_{1:t})} \left( p(m_i | z_t) = \frac{p(z_t | m_i) p(m_i)}{p(z_t)} \right)$$

Answer.  $\frac{p(z_t | m_i, x_t) p(m_i | z_{1:t-1}, x_{1:t-1})}{p(z_t | z_{1:t-1}, x_{1:t})}$

at  $x_t$ , no  $z_t$ . so  $x_t$  is not useful

• Do the same for  $\gamma_{mi}$

$$P(\gamma_{mi} | Z_{1:t}, X_{1:t}) = \frac{P(z_t | \gamma_{mi}, x_t) P(\gamma_{mi} | Z_{1:t-1}, X_{1:t-1})}{P(z_t | Z_{1:t-1}, X_{1:t-1})}$$

• ratio.

$$\frac{P(mi | Z_{1:t}, X_{1:t})}{1 - P(mi | Z_{1:t}, X_{1:t})} = \frac{\frac{P(z_t | mi, x_t) P(mi | Z_{1:t-1}, X_{1:t-1})}{P(z_t | Z_{1:t-1}, X_{1:t-1})}}{\frac{P(z_t | \gamma_{mi}, x_t) P(\gamma_{mi} | Z_{1:t-1}, X_{1:t-1})}{P(z_t | Z_{1:t-1}, X_{1:t-1})}} = \underbrace{\frac{P(z_t | mi, x_t)}{P(z_t | \gamma_{mi}, x_t)}}_{\text{bad.}} \underbrace{\frac{P(mi | Z_{1:t-1}, X_{1:t-1})}{P(\gamma_{mi} | Z_{1:t-1}, X_{1:t-1})}}_{\text{good}}$$

• Bayes for swapping:  $P(z_t | mi, x_t) \rightarrow P(mi | z_t, x_t)$ . (You see why swapping later)

$$\begin{aligned} P(z_t | mi, x_t) &= \frac{P(mi | z_t, x_t) \cdot P(z_t | x_t)}{P(mi | x_t)} \\ &= \frac{P(mi | z_t, x_t) P(z_t | x_t)}{P(mi)} \end{aligned}$$

marker: if no observation.  $x_t$  does not help.

$$\Rightarrow P(z_t | \gamma_{mi}, x_t) = \frac{P(\gamma_{mi} | z_t, x_t) P(z_t | x_t)}{P(\gamma_{mi})}$$

$$\Rightarrow \frac{P(mi | Z_{1:t}, X_{1:t})}{P(\gamma_{mi} | Z_{1:t}, X_{1:t})} = \underbrace{\frac{P(mi | z_t, x_t)}{P(\gamma_{mi} | z_t, x_t)}}_{\text{current observation and state}} \underbrace{\frac{P(\gamma_{mi})}{P(mi)}}_{\text{prior}} \underbrace{\frac{P(mi | Z_{1:t-1}, X_{1:t-1})}{P(\gamma_{mi} | Z_{1:t-1}, X_{1:t-1})}}_{\text{recursion}}$$

Recursive: for the task of  $P(mi | Z_{1:t}, X_{1:t})$ , we can keep its history. for each new state, I only need prior and ~~current~~ (obs, pos) for that timestamp to solve the mapping problem

• Ratio to probability

$$\frac{p(x)}{1-p(x)} = Y \Rightarrow p(x) = \frac{1}{1+Y}$$

$$\Rightarrow P(mi | Z_{1:t}, X_{1:t}) = \left[ 1 + \frac{1 - P(mi | z_t, x_t)}{P(mi | z_t, x_t)} \frac{P(mi)}{1 - P(mi)} \frac{1 - P(mi | Z_{1:t-1}, X_{1:t-1})}{P(mi | Z_{1:t-1}, X_{1:t-1})} \right]^{-1}$$



• For efficiency reason. We use log-odds notation.

$$l(m_i | z_{1:t}, x_{1:t}) \equiv \log \left( \frac{p(m_i | z_{1:t}, x_{1:t})}{1 - p(m_i | z_{1:t}, x_{1:t})} \right)$$

thus the transform becomes:

$$l(x) = \log \frac{p(x)}{1-p(x)} \Rightarrow p(x) = \frac{1}{1 + \exp(-l(x))}$$

And the formula becomes

$$l(m_i | z_{1:t}, x_{1:t}) \equiv l_{t,i}$$

$$= \underbrace{l(m_i | z_t, x_t)}_{\text{inverse sensor model}} + \underbrace{l(m_i | z_{1:t-1}, x_{1:t-1})}_{\text{recursive term}} - \underbrace{l(m_i)}_{\text{prior}}$$

⇒ Algorithm. given  $\{l_{t,i}\}$ ,  $x_t$ ,  $z_t$ .

for each cell  $m_i$

if  $m_i$  is in perceptum field of  $z_t$ .

$$l_{t,i} = l_{t,i} + \text{inv-sensor-model}(m_i, x_t, z_t) - l_0.$$

else

$$l_{t,i} = l_{t-1,i}.$$

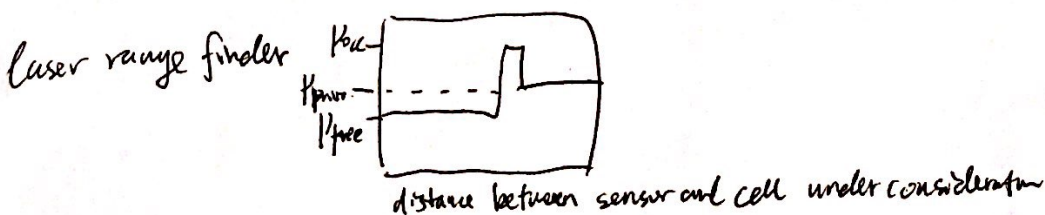
end

end.

Remark: There are many assumptions in this algorithm

- ① Grid cells are binary
- ② Grid cells are independent
- ③ Grid cells are static → we don't have a prediction step.
- ④ We know the poses perfectly → not true

4. Inverse sensor model example



## 5. Scan Match.

In reality, motion is noisy. while sensor is rather precise

=> scan matching tries to incrementally align two scans or a ~~scan~~ map to a scan without revising past map

It is in essence, pose correction.

$$x_t^* = \underset{x_t}{\operatorname{argmax}} \left\{ \underbrace{p(z_t | x_t, m_{t-1})}_{\text{measurement}} \underbrace{p(x_t | u_{t-1}, x_{t-1}^*)}_{\text{motion}} \right\}$$

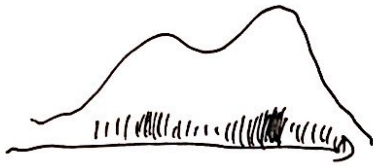
However, scan matching can only have locally consistent estimates and is not sufficient to build a large consistent map.

## Particle Filters

1. Kalman Filter Assumes a Gaussian model. Particle Filter uses a non-parametric approach for approximating a distribution.

• Natural idea would be to use samples  $\mathcal{X} = \{ \underset{\substack{\uparrow \\ \text{particle} \\ \text{sample}}}{x^{[j]}}, \underset{\substack{\uparrow \\ \text{weight. (we will see later)} \\ \text{(what this means)}}}{w^{[j]}} \} \quad j=1 \dots J.$

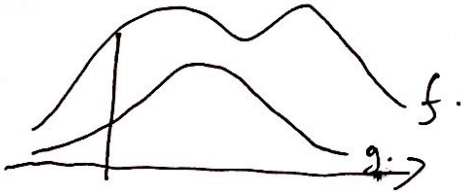
• The more particles fall into a region, the higher the probability of the region.



2. Question Comes: How to obtain the sample, in a way that makes sense?

• We know how to sample a gaussian uniformly:  $x \leftarrow \frac{1}{\sqrt{2}} \sum_{i=1}^{12} \text{rand}(-0.6, 1).$

• Importance Sampling: use a different distribution  $g$  to generate sample from  $f$ . And we know how to sample well in  $g$ .



① introduce weight:  $w = \frac{f}{g}.$

each sample will have a weight

② weights can be normalized by  $\tilde{w}_i = \frac{w_i}{\sum_i w_i}$

③ weights and samples together characterizes a distribution.  $p(x) = \sum_{j=1}^J w^{[j]} \delta_{x^{[j]}}(x).$  (weighted counts of particles in that region)

④ High weight means: one sample may "effectively" count for several samples.

~~This means we need to know our target f.~~

Obs: for <sup>high dimension distribution</sup> large variance distribution, we need many samples for a good approximation

3. ~~However we don't want~~ We can use a resampling strategy to implicitly weight the samples.

Resampling: Given  $\langle x_t^{[j]}, w_t^{[j]} \rangle \quad j=1 \dots J. = \mathcal{X}_t$

Repeat J times: draw  $i \in 1 \dots J$  with probability  $\propto w_t^{[i]}$  and add sample to  $\mathcal{X}_t$ .

Obs:



① effect of resampling is to replace unlikely samples by more likely ones "survival of fittest"

② Reduce number of samples to maintain.

⇒ particle Filter Algorithm. ( purpose: use sampling method to approximate a distribution. )  
 Assumption: have proposal. have target.

$$\bar{X}_t = X_t = \phi$$

sample  $X_t$ : for  $j=1:J$ .

sample  $X_t^{(j)} \sim \pi(x_t) \leftarrow$  proposal.

$$w_t^{(j)} = \frac{p(X_t^{(j)})}{\pi(X_t^{(j)})} \leftarrow \text{target.}$$

⇒  $\bar{X}_t = \bar{X}_t + \langle X_t^{(j)}, w_t^{(j)} \rangle$  Add to  $\bar{X}_t$ .  
 end.

② resample.

for  $j=1:J$ .

draw  $i \in 1 \dots J$  with prob  $w_t^{(i)}$   
 add sample to  $X_t$ .

end.

A well implemented algorithm

4. How to do the resampling?

Stochastic universal sampling.  
 (low variance)



5. Apply Particle Filter for Localization

Given  $X_{t-1}, u_t, z_t$ . find  $X_t$  (recursive filter)

① where do we sample from?  $X_t^{(i)} \sim p(X_t | u_t, X_{t-1}^{(i)})$  odometry model

② How to compute weight?  $w_t^{(i)} = p(z_t | X_t^{(i)})$  observation for correction

Interpretation: Given current state is  $X_t^{(i)}$ , How likely will I observe  $z_t$ ?

obs: each particle is a trajectory proposal. Then they survive according to "survival of fittest"



It's important to sample here.  
 Not translate the samples  
 (just)

③ remain filters whose observation are close to our sensor observation  
 we are here!

# Fast SLAM 1.0.

## 1. Particle Filter.

- 1.1. Three steps: Sample from proposal distribution.  
Compute importance weights  
Resampling.

1.2. Particle Filter Can be used for localization: Given  $\lambda_{t-1}, u_t, z_t$ , find  $x_t$ .

1.3. works well in low dimension, but sample a high-dim distribution is hard.

1.4. Jump to SLAM.

localization  $\rightarrow$  SLAM  
 $x_t = \begin{bmatrix} x \\ y \\ \theta \end{bmatrix}$   $x = (x_{1:t}, m_{1,x}, m_{1,y}, \dots, m_{n,x}, m_{n,y})^T$   
 high-dim state vector  
 solving this issue is important.

## 2. Rao-Blackwellization

2.1. Idea: mapping with known poses is simple (grid maps). so we can focus our state as  $x = (x_{1:t})^T$ .

This is to say, each sample represents a possible trajectory. And if this is true, we apply mapping with known poses, we get a map from each sample and compare with what we see.

## 2.2. Mathematics.

$$p(a,b) = p(b|a) p(a).$$

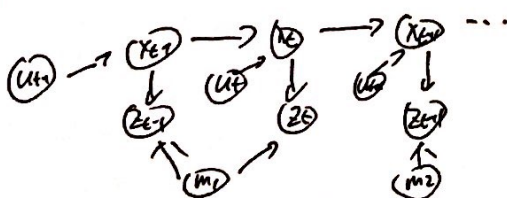
If  $p(b|a)$  can be easily computed, then sample  $p(a)$  only. And compute  $p(b|a)$  for each sample. In essence, we sampled  $p(a,b)$ .

## 2.3. RB For SLAM.

$$p(x_{0:t}, m_{1:M} | z_{1:t}, u_{1:t}) = \underbrace{p(x_{0:t} | z_{1:t}, u_{1:t})}_{\text{particle}} \underbrace{p(m_{1:M} | x_{0:t}, z_{1:t})}_{\text{mapping}}.$$

when poses are known  
controls can be ignored

## 2.4. The ~~dependence~~ model.



$x_{0:t}$  are known.  $u_{1:t}$  are independent.  
 Landmarks are also independent

## 2.5. Computer.

$$P(x_{0:t}, m_{1:M} | z_{1:t}, u_{1:t}) = \underbrace{P(x_{0:t} | z_{1:t}, u_{1:t})}_{\text{particle filter MCL.}} \prod_{i=1}^M P(m_i | x_{0:t}, z_{1:t}).$$

$\nearrow$  2x2 covariance matrix  
 $\nearrow$  2-dim EKF

build map for.

Note also each sample is a path hypothesis but we don't sense past pose. So we only need to maintain a 3-dim pose vector for the next state

## 3. Fast SLAM algorithm

### 3.1. particles.

state:  $\boxed{x, y, \theta, m_1, m_2, \dots, m_M}$

$\nearrow$  2x2 EKF.

### 3.2. Sampling:

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

for each sample, ~~use~~ <sup>apply</sup> odometry model, draw sample to get  $x_t^{(k)}$

### 3.3. Importance weight:

$$w^{(k)} = |2\pi\Omega|^{-\frac{1}{2}} \exp \left\{ -\frac{1}{2} (z_t - \hat{z}^{(k)})^T \Omega^{-1} (z_t - \hat{z}^{(k)}) \right\}$$

cannot  $\downarrow$  predicted obs  
 $\downarrow$   
 actual obs.  $\nearrow$  every particle has his own prediction.

### 3.4. The Algorithm.

Known:  $z_t, C_t, u_t, x_{t-1}$ .

#### ① Sample:

for  $k=1$  to  $N$  do

Let  $\langle x_{t-1}^{(k)}, \mu_{t-1}^{(k)}, \Sigma_{t-1}^{(k)}, \dots \rangle$  be particle  $k$  in  $\chi_{t-1}$ .

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

#### ② Detect:

$\hat{j} = C_t$  observed feature.

if feature  $j$  never seen before.



Initialize.

$$\mu_{j,t}^{[k]} = h^{-1}(z_t, x_t^{[k]})$$

$$H = h'(\mu_{j,t}^{[k]}, x_t^{[k]})$$

$$\Sigma_{j,t}^{[k]} = H^{-1} \Omega_t (H^{-1})^T$$

$$w^{[k]} = p_0.$$

else. update landmark:

$$\langle \mu_{j,t}^{[k]}, \Sigma_{j,t}^{[k]} \rangle = \text{UKF-Update}$$

$$w^{[k]} = (2\pi\Omega)^{-\frac{1}{2}} \exp \left\{ -\frac{1}{2} (z_t - \hat{z}^{[k]})^T \Omega^{-1} (z_t - \hat{z}^{[k]}) \right\}$$

$$\Omega = H \Sigma_{j,t+1}^{[k]} H^T + \Omega_t. \quad (\text{Do not update})$$

endif.

leave all unobserved features  $j'$  as they are.

endfor.

$$X_t = \text{Resample} \left( \langle x_t^{[k]}, \langle \mu_{j,t}^{[k]}, \Sigma_{j,t}^{[k]} \rangle, \dots, w^{[k]} \rangle_{k=1 \dots N} \right)$$

UKF-update

$$\hat{z}^{[k]} = h(\mu_{j,t+1}^{[k]}, x_t^{[k]}) \quad \text{--- Mousner predict}$$

$$H = h'(\mu_{j,t+1}^{[k]}, x_t^{[k]}) \quad \text{--- Jacobian}$$

$$\Omega = H \Sigma_{j,t+1}^{[k]} H^T + \Omega_t \quad \text{--- moment covariance.}$$

$$K = \Sigma_{j,t+1}^{[k]} H^T \Omega^{-1} \quad \text{--- Kalman Gain}$$

$$\left. \begin{aligned} \mu_{j,t+1}^{[k]} &= \mu_{j,t+1}^{[k]} + K(z_t - \hat{z}^{[k]}) \\ \Sigma_{j,t+1}^{[k]} &= (I - KH) \Sigma_{j,t+1}^{[k]} \end{aligned} \right\} \text{update}$$

4. Importance weights Derivation

$$\text{4.P. } w^{[k]} = \frac{\text{target}(x^{[k]})}{\text{proposal}(x^{[k]})} = \frac{P(x_{1:t} | z_{1:t}, u_{1:t})}{P(x_{1:t} | z_{1:t-1}, u_{1:t})}$$

Proposal is used step by step.

$$p(x_{1:t} | z_{1:t-1}, u_{1:t}) = \underbrace{p(x_t | x_{t-1}, u_t)}_{\text{from } x_{t-1} \text{ to } \bar{x}_t} \underbrace{p(x_{1:t-1} | z_{1:t-1}, u_{1:t-1})}_{x_{t-1}}$$

$$\Rightarrow w^{[k]} = \frac{p(x_{1:t}^{[k]} | z_{1:t}, u_{1:t})}{p(x_t^{[k]} | x_{t-1}, u_t) p(x_{1:t-1}^{[k]} | z_{1:t-1}, u_{1:t-1})}$$

$$\stackrel{\text{Bayes.}}{=} \frac{\eta p(z_t | x_{1:t}^{[k]}, z_{1:t-1}) p(x_t | x_{t-1}^{[k]}, u_t)}{p(x_t^{[k]} | x_{t-1}^{[k]}, u_t)} \frac{p(x_{1:t-1}^{[k]} | z_{1:t-1}, u_{1:t-1})}{p(x_{1:t-1}^{[k]} | z_{1:t-1}, u_{1:t-1})}$$

$$= \eta p(z_t | x_{1:t}^{[k]}, z_{1:t-1})$$

$$= \eta \int p(z_t | x_{1:t}^{[k]}, z_{1:t-1}, m_j) p(m_j | x_{1:t}^{[k]}, z_{1:t-1}) dm_j.$$

$$= \eta \int \underbrace{p(z_t | x_t^{[k]}, m_j)}_{\substack{\text{mean} \\ \mathcal{N}(z_t; \bar{z}_t^{[k]}, \Sigma_t^{[k]})}} \underbrace{p(m_j | x_{1:t-1}^{[k]}, z_{1:t-1})}_{\substack{\text{2nd GKF.} \\ \mathcal{N}(m_j; \mu_{j,t+1}^{[k]}, \Sigma_{j,t+1}^{[k]})}} dm_j.$$

$$\Rightarrow \mathcal{Q} = H \Sigma_{j,t+1}^{[k]} H^T + \mathcal{Q}_t.$$

$$\Rightarrow w^{[k]} \propto |2\pi\mathcal{Q}|^{-\frac{1}{2}} \exp\left\{-\frac{1}{2}(z_t - \bar{z}_t^{[k]})^T \mathcal{Q}^{-1} (z_t - \bar{z}_t^{[k]})\right\}$$

5. Advantages

① per-particle data association. simple but effective