Lecture 8, Oct 3, 2023

Localization

- Localization is the process of determining where the robot is
 - Do we already have a map (i.e. landmarks) or do we need to build one?
 - How do we measure uncertainty arising from sensors and actuators?
 - How do we formulate the best estimate for localization from uncertain measurements?
- Any sensor measurement will invariably be corrupted by noise to some extent
 - Measurements are often distributed according to a Gaussian, due to the central limit theorem

Propagation of Error – Odometry Example

- How does uncertainty in measurements propagate?
- The covariance matrix generalizes variance to multiple dimensions

$$\begin{array}{l} - \boldsymbol{\Sigma} = \mathbb{E}[(\boldsymbol{x} - \mathbb{E}(\boldsymbol{x}))(\boldsymbol{x} - \mathbb{E}(\boldsymbol{x}))^T] \\ = \begin{bmatrix} \sigma_{x_1}^2 & \sigma_{x_1x_2} & \cdots \\ \sigma_{x_2x_1} & \sigma_{x_2}^2 & \cdots \\ \vdots & \vdots & \ddots \end{bmatrix} \end{array}$$

- If we take the covariance between two different variables, it is known as the cross-covariance
- $\operatorname{cov}(\boldsymbol{x}, \boldsymbol{y}) = \mathbb{E}[(\boldsymbol{x} \boldsymbol{\mu}_x)(\boldsymbol{y} \boldsymbol{\mu}_y)^T]$
- Note some important properties of covariance:

1.
$$\Sigma = \mathbb{E}[xx^{T}] - \mu\mu^{T}$$

- 2. $\boldsymbol{\Sigma}^T = \boldsymbol{\Sigma} \geq 0$, i.e. the covariance matrix is semi-definite
- 3. $\operatorname{cov}(\boldsymbol{x}, \boldsymbol{y}) = \operatorname{cov}(\boldsymbol{y}, \boldsymbol{x})^T$
- 4. $\operatorname{cov}(\boldsymbol{x}_1 + \boldsymbol{x}_2, \boldsymbol{y}) = \operatorname{cov}(\boldsymbol{x}_1, \boldsymbol{y}) + \operatorname{cov}(\boldsymbol{x}_2, \boldsymbol{y})$, i.e. covariance is bilinear
- 5. $\operatorname{cov}(Ax + a, By + b) = A \operatorname{cov}(x, y)B^T$

6. $\operatorname{cov}(\boldsymbol{x}, \boldsymbol{y}) = \boldsymbol{0}$ if \boldsymbol{x} and \boldsymbol{y} are independent (but a zero covariance does not mean no correlation)

- Let $\boldsymbol{y} = \boldsymbol{f}(\boldsymbol{x})$, then in general we can see how $\boldsymbol{\Sigma}_{\boldsymbol{y}}$ relates to $\boldsymbol{\Sigma}_{\boldsymbol{x}}$
 - By Taylor expansion $\boldsymbol{y} = \boldsymbol{y}_0 + (\vec{\nabla} \boldsymbol{f})_0 (\boldsymbol{x} \boldsymbol{x}_0)$
 - * Then $(\vec{\nabla} \boldsymbol{f})_0 \boldsymbol{x} = \boldsymbol{A} \boldsymbol{x}$ and $\boldsymbol{y}_0 (\vec{\nabla} \boldsymbol{f})_0 \boldsymbol{x}_0 = \boldsymbol{a}$
 - * By property 5 above, $\Sigma_y = (\vec{\nabla} f)_0 \Sigma_x (\vec{\nabla} f)_0^T$
- Consider the problem of determining pose using only odometry, i.e. movement of the wheels $\Delta s = \begin{bmatrix} \Delta s_l \\ \Delta s_r \end{bmatrix}$

$$-\Delta \boldsymbol{x} = \begin{bmatrix} \Delta x \\ \Delta y \\ \Delta \theta \end{bmatrix} = \begin{bmatrix} \frac{\Delta s_r + \Delta s_l}{2} \cos \theta \\ \frac{\Delta s_r + \Delta s_l}{2} \sin \theta \\ \frac{\Delta s_r - \Delta s_l}{b} \end{bmatrix} = \begin{bmatrix} \Delta s \cos \theta \\ \Delta s \sin \theta \\ \frac{\Delta s_r - \Delta s_l}{b} \end{bmatrix}$$

• Our new position is given by $\Delta x' = f(x + \Delta s)$

- Linearize:
$$\mathbf{x}' = \mathbf{x} + (\vec{\nabla}_{\Delta s} \mathbf{f})_0 \Delta \mathbf{s}$$
 where $\vec{\nabla}_{\Delta s} \mathbf{f}$ is the Jacobian $\begin{bmatrix} \frac{1}{2} \cos \theta & \frac{1}{2} \cos \theta \\ \frac{1}{2} \sin \theta & \frac{1}{2} \sin \theta \\ -b^{-1} & b^{-1} \end{bmatrix}$

– Assume uncorrelated odometry error and $\boldsymbol{\Sigma}_{\Delta} = \begin{bmatrix} \sigma_{\Delta,l}^{z} & 0\\ 0 & \sigma_{\Delta,r}^{2} \end{bmatrix}$

- Error propagates as $\Sigma_{x'} = \Sigma_x + (\vec{\nabla}_{\Delta s} f) \Sigma_{\Delta} (\vec{\nabla}_{\Delta_s} f)^T$ Notice that the part we add is always positive due to the positive-semidefiniteness of the covariance, so the error always grows!

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• This means using odometry alone, our estimate of where the robot is will get worse with time

1D Kalman Filtering

- If we have n measurements for a static variable x, how do we obtain the best estimate \hat{x} ?
 - We can try to minimize $e = \sum_{k=1}^{n} w_k (\hat{x} x_k)^2$, i.e. weighted least squares
 - The weight can be $w_k = \frac{1}{\sigma_k^2}$, so that measurements with higher variance (uncertainty) are weighted less

- The solution is given by
$$\hat{x} = \frac{\sum_k \sigma_k^{-2} x_k}{\sum_k \sigma_k^{-2}}$$

- * This is a weighted average of all the x_k with weights $\frac{1}{\sigma_k^2}$
- Consider the case where we have only 2 measurements, then $\hat{x} = \frac{\sigma_2^2}{\sigma_1^2 + \sigma_2^2} x_1 + \frac{\sigma_1^2}{\sigma_1^2 + \sigma_2^2} x_2$
 - Then the variance of \hat{x} is var $\hat{x} = \left(\frac{\sigma_2^2}{\sigma_1^2 + \sigma_2^2}\right)^2 \sigma_1^2 + \left(\frac{\sigma_1^2}{\sigma_1^2 + \sigma_2^2}\right)^2 \sigma_2^2 = \frac{\sigma_1^2 \sigma_2^2}{\sigma_1^2 + \sigma_2^2}$
 - But note, this is less than both σ_1^2 and σ_2^2 !
- Note also $\hat{x} = \frac{\sigma_2^2}{\sigma_1^2 + \sigma_2^2} x_1 + \frac{\sigma_1^2}{\sigma_1^2 + \sigma_2^2} x_2 = x_1 + \frac{\sigma_1^2}{\sigma_1^2 + \sigma_2^2} (x_2 x_1)$ This is a much many conversiont form for us since makes the
 - This is a much more convenient form for us, since we've turned it from batch form (needing all measurements at once) into a recursive form (where we can continuously update)
- If we have $\hat{x}_k, \hat{\sigma}_k$ as the previous estimate at the current timestep, and we get a new measurement x_{k+1} with variance σ_{k+1}^2 then we can update:

$$- \hat{x}_{k+1} = \hat{x}_k + \frac{\hat{\sigma}_k^2}{\hat{\sigma}_k^2 + \sigma_{k+1}^2} (x_{k+1} - \hat{x}) = \hat{x}_k + W_{k+1} (x_{k+1} - \hat{x}_k) - \hat{\sigma}_{k+1} = \frac{\hat{\sigma}_k^2 \sigma_{k+1}^2}{\hat{\sigma}_k^2 + \sigma_{k+1}^2} = \hat{\sigma}_k^2 - W_{k+1} \hat{\sigma}_k^2$$

- -W is known as the Kalman gain
- We can see that this is similar to a feedback control law the correction to the state is the gain multiplied by the "error"
- Kalman filtering is a special case of Bayesian filtering, where the distribution is a Gaussian
- But we still haven't accounted for the fact that x may be dynamic, i.e. it can evolve over time; to account for this, we will predict what the new state should be based on the old estimate, and then compute the error from the measurement of the new state
 - Consider the 1D state update equation $x_{k+1} = x_k + u_k + v_k$ where u_k is the control input and v_k is some noise
 - * $v_k \sim \mathcal{N}(0, \varsigma_k^2)$, i.e. normally distributed, zero-mean with variance ς_k^2
 - * Assume that u_k can be accurately delivered, i.e. there is no noise
 - Let $\hat{x}_{k|k}$ be the estimate of x at step k, given measurements $\{x_0, x_1, \ldots, x_k\}$
 - Let $\hat{x}_{k+1|k}$ be the prediction of x at step k+1, given measurements $\{x_0, x_1, \ldots, x_k\}$ * $\hat{x}_{k+1|k} = \hat{x}_{k|k} + u_k$ (note since the noise is zero-mean, we can disregard it)
 - Now to get $\hat{x}_{k+1|k+1}$, we can use the same Kalman update formula as above

*
$$\hat{x}_{k+1|k+1} = \hat{x}_{k+1|k} + W_{k+1}(x_{k+1} - \hat{x}_{k+1|k})$$

* $\hat{\sigma}_{k+1|k+1}^2 = \hat{\sigma}_{k+1|k}^2 - W_{k+1}\hat{\sigma}_{k+1|k}^2$

• Intuitively, Kalman filters combine an estimate and a new measurement, both of which have some uncertainty, and finds the most likely new state according to the distributions of error in both

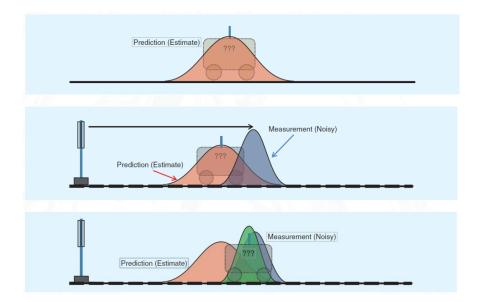


Figure 1: Diagram of Kalman filtering.