

Lecture 2: Probability

AER1513: State Estimation

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Outline

Lecture 2: Probability

- PDF Definitions

- Bayes Rule and Inference

- Gaussian PDFs

- Bayesian inference for Gaussians

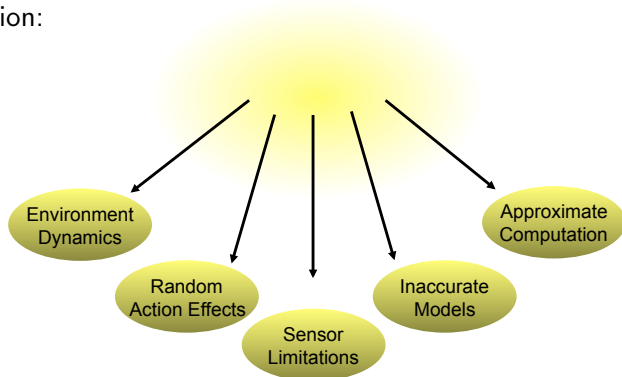
- More Gaussian tools

- Quantifying Uncertainty

- Certainty Lower Bound

It's an uncertain world

- there are several sources of uncertainty we must overcome in estimation:



- we'll need some probabilistic machinery to acknowledge and manage uncertainty

Probability densities represent uncertainty in state

- we say that a **random variable**, x , is distributed according to a particular **probability density function**
- let $p(x)$ be a probability density function (PDF) for the random variable, x , over the interval $[a, b]$
- this is a non-negative function that satisfies the **axiom of total probability**:

$$\int_a^b p(x) dx = 1 \quad (1)$$

We use PDFs to represent the likelihood of x being in all possible states in the interval, $[a, b]$.

Robot in a hallway

- this robot has a prior map of the hallway (i.e., knows where the doors are) and then detects a door

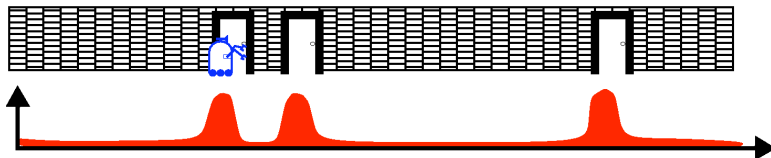
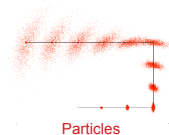
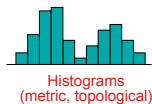
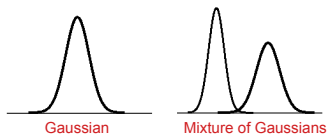


image: Thrun et al. (2006)

- it will have a PDF, $p(x)$, with three peaks representing that it is likely that it is near a door

Representations

- as we'll see, we can't represent PDFs perfectly in a computer so we need to choose an approximation

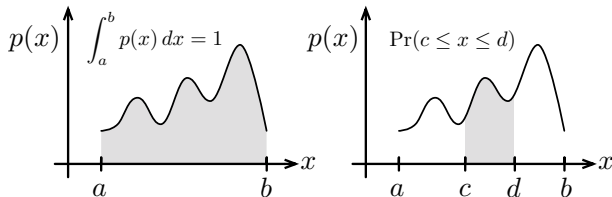


- each has its pros and cons, depending on the application

Probability is area under the curve

- note that **probability density** is not **probability**
- **probability** is given by the area under the density function
- for example, the probability that x lies between c and d , $\Pr(c \leq x \leq d)$, is given by

$$\Pr(c \leq x \leq d) = \int_c^d p(x) dx \quad (2)$$



Cumulative density and quantile functions

- the **cumulative density function (CDF)**, $P(x)$, is given by

$$P(x) = \Pr(x' \leq x) = \int_{-\infty}^x p(x') dx' \quad (3)$$

which is the area under the density function, $p(x)$, up to a particular x

- the **quantile function**, $Q(y)$, is the inverse (if it exists) of the CDF:

$$Q(y) = P^{-1}(y) \quad (4)$$

where the range is between 0 and 1; we will use this later to sample from a distribution and also to perform statistical hypothesis tests

There are often strings attached

- we can introduce a **conditioning variable** to our PDFs
- let $p(x|y)$ be a PDF over $x \in [a, b]$ conditioned on $y \in [r, s]$ such that

$$(\forall y) \quad \int_a^b p(x|y) dx = 1 \quad (5)$$

- this tells us about the likelihood of x given a particular value of y
- for example, x could be a robot position and y could be some sensor readings

The curse of dimensionality

- we may also denote **joint probability densities** for N -dimensional continuous variables in our framework as $p(\mathbf{x})$ where

$$\mathbf{x} = (x_1, \dots, x_N) \quad (6)$$

with $x_i \in [a_i, b_i]$

- note that we can also use the notation

$$p(x_1, x_2, \dots, x_N) \quad (7)$$

in place of $p(\mathbf{x})$

- sometimes we even mix and match the two and write

$$p(\mathbf{x}, \mathbf{y}) \quad (8)$$

for the joint density of \mathbf{x} and \mathbf{y}

Dimensions don't trump axioms

- in the N -dimensional case, the **axiom of total probability** requires

$$\begin{aligned} \int_{\mathbf{a}}^{\mathbf{b}} p(\mathbf{x}) d\mathbf{x} \\ = \int_{a_N}^{b_N} \cdots \int_{a_2}^{b_2} \int_{a_1}^{b_1} p(x_1, x_2, \dots, x_N) dx_1 dx_2 \cdots dx_N = 1 \quad (9) \end{aligned}$$

where $\mathbf{a} = (a_1, \dots, a_N)$ and $\mathbf{b} = (b_1, \dots, b_N)$

Factoring a PDF

- we can always **factor a joint probability density** into a conditional and an unconditional factor:

$$p(\mathbf{x}, \mathbf{y}) = p(\mathbf{x}|\mathbf{y})p(\mathbf{y}) = p(\mathbf{y}|\mathbf{x})p(\mathbf{x}) \quad (10)$$

- in the specific case that \mathbf{x} and \mathbf{y} are **statistically independent**, we can factor the joint density as follows:

$$p(\mathbf{x}, \mathbf{y}) = p(\mathbf{x})p(\mathbf{y}) \quad (11)$$

or put another way, $p(\mathbf{x}|\mathbf{y}) = p(\mathbf{x})$ and $p(\mathbf{y}|\mathbf{x}) = p(\mathbf{y})$

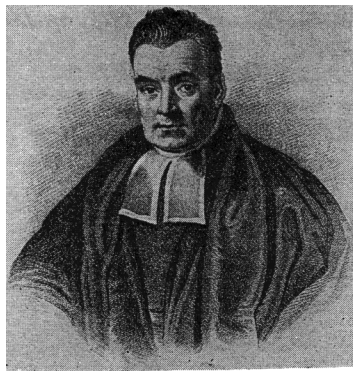
Bayes' rule

- restating the factored expression:

$$p(\mathbf{x}, \mathbf{y}) = p(\mathbf{x}|\mathbf{y})p(\mathbf{y}) = p(\mathbf{y}|\mathbf{x})p(\mathbf{x}) \quad (12)$$

- **Bayes' rule** (Bayes, 1764) follows by rearranging:

$$p(\mathbf{x}|\mathbf{y}) = \frac{p(\mathbf{y}|\mathbf{x})p(\mathbf{x})}{p(\mathbf{y})} \quad (13)$$



REV. T. BAYES

Bayesian inference

- we use **Bayes' rule** to **infer** one probability density function from another:

$$p(\mathbf{x}|\mathbf{y}) = \frac{p(\mathbf{y}|\mathbf{x})p(\mathbf{x})}{p(\mathbf{y})} \quad (14)$$

- $p(\mathbf{x})$ is called the **prior density** (does not incorporate any data)
- $p(\mathbf{x}|\mathbf{y})$ is called the **posterior density** (incorporates data)
- $p(\mathbf{y}|\mathbf{x})$ is a **generative model** (e.g., sensor model)
- $p(\mathbf{y})$, the denominator, is discussed on the next slide

Bayesian inference is the cornerstone of modern state estimation and machine learning.

On the margins

- we can compute the denominator in Bayes' rule using the idea of **marginalization** of a joint PDF,

$$p(\mathbf{x}, \mathbf{y}) \tag{15}$$

- if we integrate over all possible values of one of the joint variables, we can compute the density over only the other:

$$\int_a^b p(\mathbf{x}, \mathbf{y}) d\mathbf{x} = \int_a^b p(\mathbf{x}|\mathbf{y})p(\mathbf{y}) d\mathbf{x} = p(\mathbf{y}) \underbrace{\int_a^b p(\mathbf{x}|\mathbf{y}) d\mathbf{x}}_1 = p(\mathbf{y}) \tag{16}$$

- this is typically the most expensive step in Bayesian estimation

Just a moment

- when working with mass distributions (a.k.a., density functions) in dynamics, we often keep track of only a few properties called the **moments** of mass (e.g., mass, center of mass, inertia matrix)
- the same is true with probability density functions
- the **zeroeth probability moment** is always 1 owing to the axiom of total probability
- the **first probability moment** is known as the **mean**, μ :

$$\mu = E[x] = \int_a^b x p(x) dx \quad (17)$$

where $E[\cdot]$ denotes the **expectation operator**

To infinity and beyond!

- the **second probability moment** is known as the **covariance matrix**, Σ :

$$\Sigma = E [(\mathbf{x} - \boldsymbol{\mu})(\mathbf{x} - \boldsymbol{\mu})^T] = \int_{\mathbf{a}}^{\mathbf{b}} (\mathbf{x} - \boldsymbol{\mu})(\mathbf{x} - \boldsymbol{\mu})^T p(\mathbf{x}) d\mathbf{x} \quad (18)$$

- in $N = 1$ dimension, this is the familiar **variance**, σ^2 , with σ the **standard deviation**
- the next two moments are called the **skewness** and **kurtosis**, but for the multivariate case, these get quite complicated and require tensor representations
- we will not need them here, but it should be mentioned that there are an infinite number of these probability moments

Sample mean and covariance

- we can draw **samples** (or **realizations**) from a PDF, which we denote as $\mathbf{x}_{\text{meas}} \leftarrow p(\mathbf{x})$; aside: how would we do this?
- to estimate the mean and covariance of random variable, \mathbf{x} , from N samples we use the **sample mean** and **sample covariance**:

$$\boldsymbol{\mu}_{\text{meas}} = \frac{1}{N} \sum_{i=1}^N \mathbf{x}_{i,\text{meas}} \quad (19a)$$

$$\boldsymbol{\Sigma}_{\text{meas}} = \frac{1}{N-1} \sum_{i=1}^N (\mathbf{x}_{i,\text{meas}} - \boldsymbol{\mu}_{\text{meas}}) (\mathbf{x}_{i,\text{meas}} - \boldsymbol{\mu}_{\text{meas}})^T \quad (19b)$$

- the sample covariance uses $N - 1$ rather than N in the denominator since it uses the sample mean, which is computed from the same measurements, resulting in a slight correlation

Statistically independent vs. uncorrelated

- two random variables, \mathbf{x} and \mathbf{y} , are **statistically independent** if their joint density factors as follows:

$$p(\mathbf{x}, \mathbf{y}) = p(\mathbf{x}) p(\mathbf{y}) \quad (20)$$

- two random variables are **uncorrelated** if

$$E[\mathbf{x}\mathbf{y}^T] = E[\mathbf{x}] E[\mathbf{y}]^T \quad (21)$$

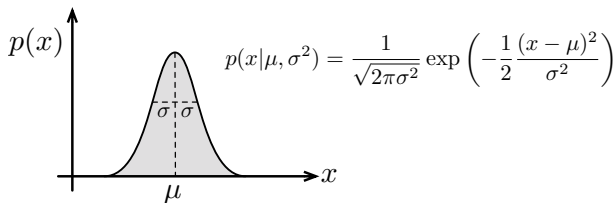
- statistically independent always implies uncorrelated, but the reverse is not always true (but sometimes it is)

Gauss' namesake

- we'll be working with **Gaussian** probability density functions
- in one dimension, a Gaussian PDF is given by

$$p(x|\mu, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{1}{2} \frac{(x - \mu)^2}{\sigma^2}\right) \quad (22)$$

where μ is the **mean** and σ^2 is the **variance** (σ is called the **standard deviation**)



Multivariate Gaussian

- a **multivariate Gaussian** probability density function, $p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma})$, over the random variable, $\mathbf{x} \in \mathbb{R}^N$, may be expressed as

$$p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) = \frac{1}{\sqrt{(2\pi)^N \det \boldsymbol{\Sigma}}} \exp \left(-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu}) \right) \quad (23)$$

- $\boldsymbol{\mu} \in \mathbb{R}^N$ is the **mean**
- $\boldsymbol{\Sigma} \in \mathbb{R}^{N \times N}$ is the (symmetric positive-definite) **covariance matrix**

Gaussian moments

- for a Gaussian we must therefore have that

$$\begin{aligned}\boldsymbol{\mu} = E[\mathbf{x}] &= \int_{-\infty}^{\infty} \mathbf{x} \frac{1}{\sqrt{(2\pi)^N \det \boldsymbol{\Sigma}}} \\ &\quad \times \exp\left(-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu})\right) d\mathbf{x} \quad (24)\end{aligned}$$

and

$$\begin{aligned}\boldsymbol{\Sigma} &= E[(\mathbf{x} - \boldsymbol{\mu})(\mathbf{x} - \boldsymbol{\mu})^T] \\ &= \int_{-\infty}^{\infty} (\mathbf{x} - \boldsymbol{\mu})(\mathbf{x} - \boldsymbol{\mu})^T \frac{1}{\sqrt{(2\pi)^N \det \boldsymbol{\Sigma}}} \\ &\quad \times \exp\left(-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu})\right) d\mathbf{x} \quad (25)\end{aligned}$$

- check for yourself!

Gaussians are the new normal

- we may also write that \mathbf{x} is **normally** (a.k.a., Gaussian) distributed using the following notation:

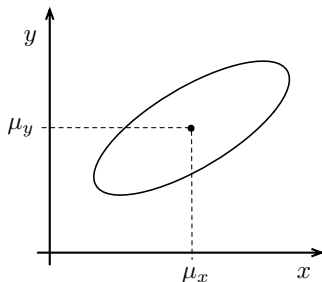
$$\mathbf{x} \sim \mathcal{N}(\boldsymbol{\mu}, \boldsymbol{\Sigma}) \quad (26)$$

- we say a random variable is **standard normally** distributed if

$$\mathbf{x} \sim \mathcal{N}(\mathbf{0}, \mathbf{1}) \quad (27)$$

where $\mathbf{1}$ is an $N \times N$ identity matrix

Joint Gaussians



- we can also have a **joint Gaussian** over a pair of variables, (\mathbf{x}, \mathbf{y}) , which we write as

$$p(\mathbf{x}, \mathbf{y}) = \mathcal{N} \left(\begin{bmatrix} \mu_x \\ \mu_y \end{bmatrix}, \begin{bmatrix} \Sigma_{xx} & \Sigma_{xy} \\ \Sigma_{yx} & \Sigma_{yy} \end{bmatrix} \right) \quad (28)$$

with $\Sigma_{yx} = \Sigma_{xy}^T$

Statistically independent vs. uncorrelated

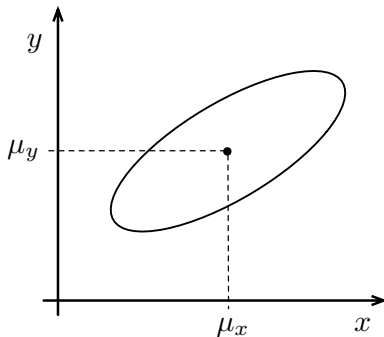
- for joint Gaussians, **statistically independent** implies **uncorrelated** and vice versa
- independence always implies uncorrelated (for any PDF)
- to go the other way, assume uncorrelated

$$\Sigma_{xy} = \Sigma_{yx}^T = \mathbf{0} \quad (29)$$

which implies independence:

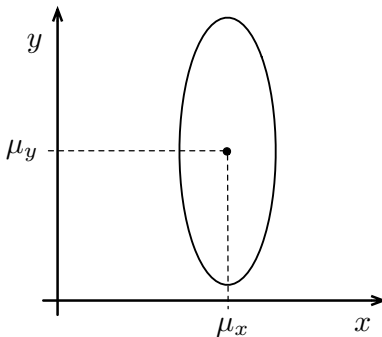
$$\begin{aligned} p(\mathbf{x}|\mathbf{y}) &= \mathcal{N}(\boldsymbol{\mu}_x + \boldsymbol{\Sigma}_{xy}\boldsymbol{\Sigma}_{yy}^{-1}(\mathbf{y} - \boldsymbol{\mu}_y), \boldsymbol{\Sigma}_{xx} - \boldsymbol{\Sigma}_{xy}\boldsymbol{\Sigma}_{yy}^{-1}\boldsymbol{\Sigma}_{yx}) \\ &= \mathcal{N}(\boldsymbol{\mu}_x, \boldsymbol{\Sigma}_{xx}) = p(\mathbf{x}) \end{aligned} \quad (30)$$

Statistically independent vs. uncorrelated



correlated

$$E[(x - \mu_x)(y - \mu_y)] \neq 0$$



uncorrelated

$$E[(x - \mu_x)(y - \mu_y)] = 0$$

Gaussian inference

- recall that we can **factor** any joint PDF according to

$$p(\mathbf{x}, \mathbf{y}) = p(\mathbf{x}|\mathbf{y}) p(\mathbf{y}) \quad (31)$$

- in the case of a joint Gaussian, the factors are

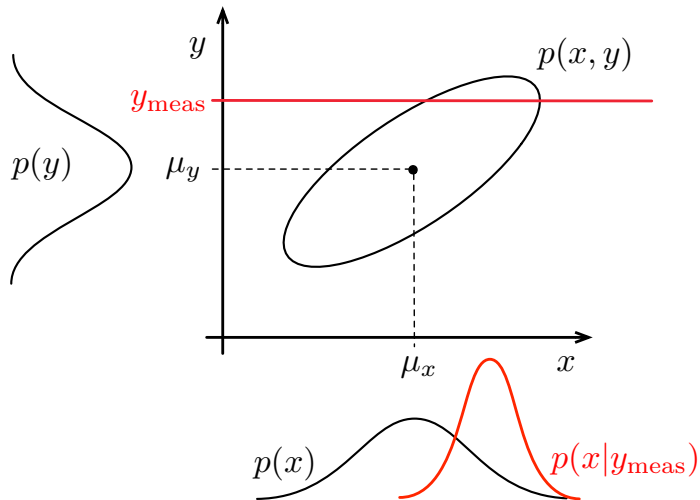
$$p(\mathbf{x}, \mathbf{y}) = \mathcal{N} \left(\begin{bmatrix} \boldsymbol{\mu}_x \\ \boldsymbol{\mu}_y \end{bmatrix}, \begin{bmatrix} \boldsymbol{\Sigma}_{xx} & \boldsymbol{\Sigma}_{xy} \\ \boldsymbol{\Sigma}_{yx} & \boldsymbol{\Sigma}_{yy} \end{bmatrix} \right) \quad (32a)$$

$$p(\mathbf{x}|\mathbf{y}) = \mathcal{N} \left(\boldsymbol{\mu}_x + \boldsymbol{\Sigma}_{xy} \boldsymbol{\Sigma}_{yy}^{-1} (\mathbf{y} - \boldsymbol{\mu}_y), \boldsymbol{\Sigma}_{xx} - \boldsymbol{\Sigma}_{xy} \boldsymbol{\Sigma}_{yy}^{-1} \boldsymbol{\Sigma}_{yx} \right) \quad (32b)$$

$$p(\mathbf{y}) = \mathcal{N} (\boldsymbol{\mu}_y, \boldsymbol{\Sigma}_{yy}) \quad (32c)$$

The expression for $p(\mathbf{x}|\mathbf{y})$ constitutes **Bayesian inference** for the case of Gaussian PDFs; we'll use this a lot!

Gaussian inference



How can you be Schur?

- to factor the joint Gaussian, we use the **Schur complement**:

$$\begin{bmatrix} \Sigma_{xx} & \Sigma_{xy} \\ \Sigma_{yx} & \Sigma_{yy} \end{bmatrix} = \begin{bmatrix} 1 & \Sigma_{xy} \Sigma_{yy}^{-1} \\ 0 & 1 \end{bmatrix} \times \begin{bmatrix} \Sigma_{xx} - \Sigma_{xy} \Sigma_{yy}^{-1} \Sigma_{yx} & 0 \\ 0 & \Sigma_{yy} \end{bmatrix} \begin{bmatrix} 1 & 0 \\ \Sigma_{yy}^{-1} \Sigma_{yx} & 1 \end{bmatrix} \quad (33)$$

- inverting this we have

$$\begin{bmatrix} \Sigma_{xx} & \Sigma_{xy} \\ \Sigma_{yx} & \Sigma_{yy} \end{bmatrix}^{-1} = \begin{bmatrix} 1 & 0 \\ -\Sigma_{yy}^{-1} \Sigma_{yx} & 1 \end{bmatrix} \times \begin{bmatrix} (\Sigma_{xx} - \Sigma_{xy} \Sigma_{yy}^{-1} \Sigma_{yx})^{-1} & 0 \\ 0 & \Sigma_{yy}^{-1} \end{bmatrix} \begin{bmatrix} 1 & -\Sigma_{xy} \Sigma_{yy}^{-1} \\ 0 & 1 \end{bmatrix} \quad (34)$$

That's for Schur

- plugging this into the quadratic part of the Gaussian PDF we have

$$\begin{aligned}\log p(\mathbf{x}, \mathbf{y}) &\propto \left(\begin{bmatrix} \mathbf{x} \\ \mathbf{y} \end{bmatrix} - \begin{bmatrix} \boldsymbol{\mu}_x \\ \boldsymbol{\mu}_y \end{bmatrix} \right)^T \begin{bmatrix} \boldsymbol{\Sigma}_{xx} & \boldsymbol{\Sigma}_{xy} \\ \boldsymbol{\Sigma}_{yx} & \boldsymbol{\Sigma}_{yy} \end{bmatrix}^{-1} \left(\begin{bmatrix} \mathbf{x} \\ \mathbf{y} \end{bmatrix} - \begin{bmatrix} \boldsymbol{\mu}_x \\ \boldsymbol{\mu}_y \end{bmatrix} \right) \\ &= \underbrace{\left(\mathbf{x} - \boldsymbol{\mu}_x - \boldsymbol{\Sigma}_{xy} \boldsymbol{\Sigma}_{yy}^{-1} (\mathbf{y} - \boldsymbol{\mu}_y) \right)^T \left(\boldsymbol{\Sigma}_{xx} - \boldsymbol{\Sigma}_{xy} \boldsymbol{\Sigma}_{yy}^{-1} \boldsymbol{\Sigma}_{yx} \right)^{-1} \left(\mathbf{x} - \boldsymbol{\mu}_x - \boldsymbol{\Sigma}_{xy} \boldsymbol{\Sigma}_{yy}^{-1} (\mathbf{y} - \boldsymbol{\mu}_y) \right)}_{\propto \log p(\mathbf{x}|\mathbf{y})} \\ &\quad + \underbrace{\left(\mathbf{y} - \boldsymbol{\mu}_y \right)^T \boldsymbol{\Sigma}_{yy}^{-1} \left(\mathbf{y} - \boldsymbol{\mu}_y \right)}_{\propto \log p(\mathbf{y})} \quad (35)\end{aligned}$$

- recall that $\log(ab) = \log(a) + \log(b)$, for a and b scalar

Normalized product of Gaussians

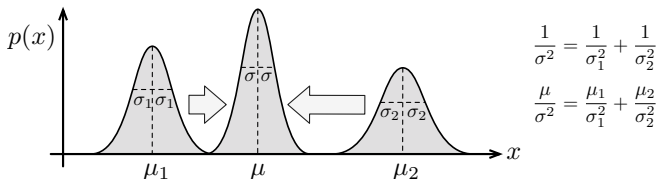
- the direct product of K Gaussian PDFs is also a Gaussian PDF:

$$p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) \equiv \prod_{k=1}^K p(\mathbf{x}|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k) \quad (36)$$

where

$$\boldsymbol{\Sigma}^{-1} = \sum_{k=1}^K \boldsymbol{\Sigma}_k^{-1}, \quad \boldsymbol{\Sigma}^{-1}\boldsymbol{\mu} = \sum_{k=1}^K \boldsymbol{\Sigma}_k^{-1}\boldsymbol{\mu}_k \quad (37)$$

- this gets used frequently in **information fusion**



Normalized product variation

– we also have that

$$\begin{aligned} & \exp \left(-\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) \right) \\ & \equiv \eta \prod_{k=1}^K \exp \left(-\frac{1}{2} (\mathbf{G}_k \mathbf{x} - \boldsymbol{\mu}_k)^T \boldsymbol{\Sigma}_k^{-1} (\mathbf{G}_k \mathbf{x} - \boldsymbol{\mu}_k) \right), \end{aligned} \quad (38)$$

where

$$\boldsymbol{\Sigma}^{-1} = \sum_{k=1}^K \mathbf{G}_k^T \boldsymbol{\Sigma}_k^{-1} \mathbf{G}_k, \quad \boldsymbol{\Sigma}^{-1} \boldsymbol{\mu} = \sum_{k=1}^K \mathbf{G}_k^T \boldsymbol{\Sigma}_k^{-1} \boldsymbol{\mu}_k, \quad (39)$$

in the case that the matrices, $\mathbf{G}_k \in \mathbb{R}^{M_k \times N}$, are present, with $M_k \leq N$; η is a normalization constant

Gaussian transformation

- we now examine **Gaussian transformation**, namely computing

$$p(\mathbf{y}) = \int_{\mathbf{x}} p(\mathbf{y}|\mathbf{x})p(\mathbf{x})d\mathbf{x}$$

where we have that

$$\begin{aligned} p(\mathbf{y}|\mathbf{x}) &= \mathcal{N}(\mathbf{g}(\mathbf{x}), \mathbf{R}) \\ p(\mathbf{x}) &= \mathcal{N}(\boldsymbol{\mu}_x, \boldsymbol{\Sigma}_{xx}) \end{aligned}$$

and $\mathbf{g}(\cdot)$ is a nonlinear map: $\mathbf{g} : \mathbf{x} \mapsto \mathbf{y}$

- this is used, for example, in the denominator when carrying out full Bayesian inference using Gaussians
- the problem is that we can't compute this integral in general, so we need to approximate it

Transformation via linearization

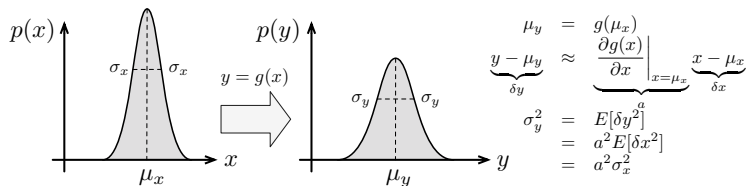
- we linearize the map such that

$$\mathbf{g}(\mathbf{x}) \approx \boldsymbol{\mu}_y + \mathbf{G}(\mathbf{x} - \boldsymbol{\mu}_x) \quad (40a)$$

$$\mathbf{G} = \left. \frac{\partial \mathbf{g}(\mathbf{x})}{\partial \mathbf{x}} \right|_{\mathbf{x}=\boldsymbol{\mu}_x} \quad (40b)$$

$$\boldsymbol{\mu}_y = \mathbf{g}(\boldsymbol{\mu}_x) \quad (40c)$$

where \mathbf{G} is the Jacobian of \mathbf{g} , with respect to \mathbf{x}



Transformation via linearization

- we have that

$$p(\mathbf{y}) = \int_{-\infty}^{\infty} p(\mathbf{y}|\mathbf{x})p(\mathbf{x})d\mathbf{x} \quad (41)$$

$$\begin{aligned} &= \eta \int_{-\infty}^{\infty} \exp \left(-\frac{1}{2} (\mathbf{y} - (\boldsymbol{\mu}_y + \mathbf{G}(\mathbf{x} - \boldsymbol{\mu}_x)))^T \right. \\ &\quad \times \left. \mathbf{R}^{-1} (\mathbf{y} - (\boldsymbol{\mu}_y + \mathbf{G}(\mathbf{x} - \boldsymbol{\mu}_x))) \right) \end{aligned} \quad (42)$$

$$\times \exp \left(-\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu}_x)^T \boldsymbol{\Sigma}_{xx}^{-1} (\mathbf{x} - \boldsymbol{\mu}_x) \right) d\mathbf{x} \quad (43)$$

$$= \rho \exp \left(-\frac{1}{2} (\mathbf{y} - \boldsymbol{\mu}_y)^T (\mathbf{R} + \mathbf{G}\boldsymbol{\Sigma}_{xx}\mathbf{G}^T)^{-1} (\mathbf{y} - \boldsymbol{\mu}_y) \right) \quad (44)$$

with ρ a normalization constant

- this is exactly Gaussian in \mathbf{y} :

$$\mathbf{y} \sim \mathcal{N}(\boldsymbol{\mu}_y, \boldsymbol{\Sigma}_{yy}) = \mathcal{N}(\mathbf{g}(\boldsymbol{\mu}_x), \mathbf{R} + \mathbf{G}\boldsymbol{\Sigma}_{xx}\mathbf{G}^T) \quad (45)$$

Sherman-Morrison-Woodbury identity

- the full derivation from the previous slide makes use of the **Sherman-Morrison-Woodbury (SMW) identity**, also sometimes called the **matrix inversion lemma**
- these identities are used frequently when manipulating expressions involving the covariance matrices associated with Gaussian PDFs:

$$(\mathbf{A}^{-1} + \mathbf{B}\mathbf{D}^{-1}\mathbf{C})^{-1} \equiv \mathbf{A} - \mathbf{A}\mathbf{B}(\mathbf{D} + \mathbf{C}\mathbf{A}\mathbf{B})^{-1}\mathbf{C}\mathbf{A} \quad (46a)$$

$$(\mathbf{D} + \mathbf{C}\mathbf{A}\mathbf{B})^{-1} \equiv \mathbf{D}^{-1} - \mathbf{D}^{-1}\mathbf{C}(\mathbf{A}^{-1} + \mathbf{B}\mathbf{D}^{-1}\mathbf{C})^{-1}\mathbf{B}\mathbf{D}^{-1} \quad (46b)$$

$$\mathbf{A}\mathbf{B}(\mathbf{D} + \mathbf{C}\mathbf{A}\mathbf{B})^{-1} \equiv (\mathbf{A}^{-1} + \mathbf{B}\mathbf{D}^{-1}\mathbf{C})^{-1}\mathbf{B}\mathbf{D}^{-1} \quad (46c)$$

$$(\mathbf{D} + \mathbf{C}\mathbf{A}\mathbf{B})^{-1}\mathbf{C}\mathbf{A} \equiv \mathbf{D}^{-1}\mathbf{C}(\mathbf{A}^{-1} + \mathbf{B}\mathbf{D}^{-1}\mathbf{C})^{-1} \quad (46d)$$

- they were originally discovered when attempting to update the inverse of a big matrix with a few entries slightly changed

Quantifying uncertainty

- often in problems of estimation, we have estimated a PDF for some random variable and then want to quantify how certain we are in, for example, the mean of that PDF
- one method of doing this is to look at the **negative entropy** or **Shannon information**, H , which is given by

$$H(\mathbf{x}) = -E[\ln p(\mathbf{x})] = -\int_a^b p(\mathbf{x}) \ln p(\mathbf{x}) d\mathbf{x}$$

- we'll make this expression specific to Gaussian PDFs next

Gaussian information

- for a Gaussian PDF, we have for the Shannon information:

$$H(\mathbf{x}) = - \int_{-\infty}^{\infty} p(\mathbf{x}) \ln p(\mathbf{x}) d\mathbf{x} \quad (47a)$$

$$= - \int_{-\infty}^{\infty} p(\mathbf{x}) \left(-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu}) - \ln \sqrt{(2\pi)^N \det \boldsymbol{\Sigma}} \right) d\mathbf{x} \quad (47b)$$

$$= \frac{1}{2} \ln ((2\pi)^N \det \boldsymbol{\Sigma}) + \int_{-\infty}^{\infty} \frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu}) p(\mathbf{x}) d\mathbf{x} \quad (47c)$$

$$= \frac{1}{2} \ln ((2\pi)^N \det \boldsymbol{\Sigma}) + \frac{1}{2} E [(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu})] \quad (47d)$$

- this term is exactly a squared **Mahalanobis distance**, which is like a squared Euclidean distance, but weighted in the middle by the inverse covariance matrix

Information manipulation

- a nice property of this quadratic function inside the expectation allows to rewrite it using the (linear) trace operator from linear algebra:

$$(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) = \text{tr} \left(\boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) (\mathbf{x} - \boldsymbol{\mu})^T \right) \quad (48)$$

- since the expectation is also a linear operator, we may interchange the order of the expectation and trace arriving at:

$$\begin{aligned} E \left[(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) \right] &= \text{tr} \left(E \left[\boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) (\mathbf{x} - \boldsymbol{\mu})^T \right] \right) \\ &= \text{tr} \left(\boldsymbol{\Sigma}^{-1} \underbrace{E \left[(\mathbf{x} - \boldsymbol{\mu}) (\mathbf{x} - \boldsymbol{\mu})^T \right]}_{\boldsymbol{\Sigma}} \right) = \text{tr} \left(\boldsymbol{\Sigma}^{-1} \boldsymbol{\Sigma} \right) = \text{tr} \mathbf{1} = N \end{aligned}$$

which is just the dimension of the variable!

The bottom line

- finally, for our Shannon information expression we have

$$H(\mathbf{x}) = \frac{1}{2} \ln((2\pi)^N \det \Sigma) + \frac{1}{2} E[(\mathbf{x} - \boldsymbol{\mu})^T \Sigma^{-1} (\mathbf{x} - \boldsymbol{\mu})] \quad (49a)$$

$$= \frac{1}{2} \ln((2\pi)^N \det \Sigma) + \frac{1}{2} N \quad (49b)$$

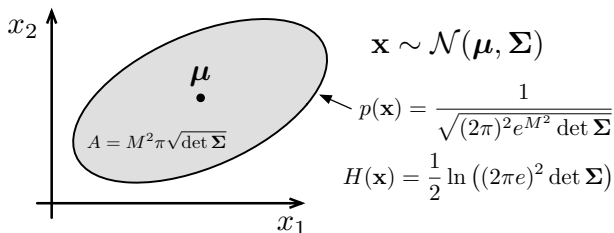
$$= \frac{1}{2} (\ln((2\pi)^N \det \Sigma) + N \ln e) \quad (49c)$$

$$= \frac{1}{2} \ln((2\pi e)^N \det \Sigma), \quad (49d)$$

which is purely a function of Σ , the covariance matrix of the Gaussian PDF

Uncertainty ellipsoid

- geometrically, we may interpret $\sqrt{\det \Sigma}$ as the volume of the **uncertainty ellipsoid** formed by the Gaussian PDF



- the geometric area inside the ellipse is

$$A = M^2 \pi \sqrt{\det \Sigma} \quad (50)$$

Equilikely contours

- note that along the boundary of the uncertainty ellipse, $p(\mathbf{x})$ is constant
- to see this, consider that the points along this ellipse must satisfy

$$(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) = M^2 \quad (51)$$

where M is a factor applied to scale the nominal ($M = 1$) covariance

- in this case, we have that

$$p(\mathbf{x}) = \frac{1}{\sqrt{(2\pi)^N e^{M^2} \det \boldsymbol{\Sigma}}} \quad (52)$$

which is independent of \mathbf{x} and thus constant

Drawing samples

- suppose we have a deterministic parameter, θ , that influences the outcome of a random variable, \mathbf{x}
- this can be captured by writing the PDF for \mathbf{x} as depending on θ :

$$p(\mathbf{x}|\theta) \tag{53}$$

- further suppose we now draw a sample, \mathbf{x}_{meas} , from $p(\mathbf{x}|\theta)$:

$$\mathbf{x}_{\text{meas}} \leftarrow p(\mathbf{x}|\theta) \tag{54}$$

- the \mathbf{x}_{meas} is sometimes called a **realization** of the random variable \mathbf{x} ; we can think of it as a ‘measurement’

As sure as sure can be

- the **Cramér-Rao lower bound (CRLB)** says that the covariance of any **unbiased estimate**, $\hat{\boldsymbol{\theta}}$ (based on the measurement, \mathbf{x}_{meas}), of the deterministic parameter, $\boldsymbol{\theta}$, is bounded by the **Fisher information matrix**, $\mathcal{I}_{\boldsymbol{\theta}}$:

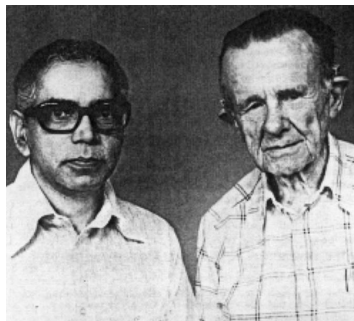
$$\text{cov}(\hat{\boldsymbol{\theta}}|\mathbf{x}_{\text{meas}}) = E \left[(\hat{\boldsymbol{\theta}} - \boldsymbol{\theta})(\hat{\boldsymbol{\theta}} - \boldsymbol{\theta})^T \right] \geq \mathcal{I}_{\boldsymbol{\theta}}^{-1}$$

where ‘unbiased’ implies $E \left[\hat{\boldsymbol{\theta}} - \boldsymbol{\theta} \right] = \mathbf{0}$ and ‘bounded’ means $\text{cov}(\hat{\boldsymbol{\theta}}|\mathbf{x}_{\text{meas}}) - \mathcal{I}_{\boldsymbol{\theta}}^{-1}$ is positive-semi-definite

- the Fisher information matrix is given by

$$\mathcal{I}_{\boldsymbol{\theta}} = E \left[\frac{\partial^2 (-\ln p(\mathbf{x}|\boldsymbol{\theta}))}{\partial \boldsymbol{\theta}^T \partial \boldsymbol{\theta}} \right]$$

What on Earth does that mean?



C. R. Rao, Harald Cramér, 1978

The **Cramér-Rao lower bound** says there is a fundamental limit on how certain we can be about an estimate of a parameter, given our measurements, regardless of the form of the estimator we use.

Our first estimation problem

- suppose that we have K samples (i.e., measurements), $\mathbf{x}_{\text{meas},k} \in \mathbb{R}^N$, drawn from a Gaussian PDF
- the K **statistically independent** random variables associated with these measurements are thus

$$(\forall k) \quad \mathbf{x}_k \sim \mathcal{N}(\boldsymbol{\mu}, \boldsymbol{\Sigma}) \quad (55)$$

- the term statistically independent implies that $E[(\mathbf{x}_k - \bar{\mathbf{x}})(\mathbf{x}_l - \bar{\mathbf{x}})^T] = \mathbf{0}$ for $k \neq l$
- now suppose our goal is to estimate the mean of this probability density function, $\boldsymbol{\mu}$, from the measurements, $\mathbf{x}_{\text{meas},1}, \dots, \mathbf{x}_{\text{meas},K}$

Joint density

- for the joint density of all the random variables, $\mathbf{x} = (\mathbf{x}_1, \dots, \mathbf{x}_K)$, we have

$$\ln p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) = -\frac{1}{2}(\mathbf{x} - \mathbf{A}\boldsymbol{\mu})^T \mathbf{B}^{-1}(\mathbf{x} - \mathbf{A}\boldsymbol{\mu}) - \ln \sqrt{(2\pi)^{NK} \det \mathbf{B}} \quad (56)$$

where

$$\mathbf{A} = \underbrace{\begin{bmatrix} \mathbf{1} & \mathbf{1} & \dots & \mathbf{1} \end{bmatrix}^T}_{K \text{ blocks}}, \quad \mathbf{B} = \text{diag} \left(\underbrace{\boldsymbol{\Sigma}, \boldsymbol{\Sigma}, \dots, \boldsymbol{\Sigma}}_{K \text{ blocks}} \right) \quad (57)$$

- in this case, we have

$$\frac{\partial^2 (-\ln p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}))}{\partial \boldsymbol{\mu}^T \partial \boldsymbol{\mu}} = \mathbf{A}^T \mathbf{B}^{-1} \mathbf{A} \quad (58)$$

Fisher information matrix

- the Fisher information matrix is

$$\mathcal{I}_{\mu} = E \left[\frac{\partial^2 (-\ln p(\mathbf{x}|\mu))}{\partial \mu^T \partial \mu} \right] \quad (59a)$$

$$= \mathbf{A}^T \mathbf{B}^{-1} \mathbf{A} \quad (59b)$$

$$= K \Sigma^{-1} \quad (59c)$$

which we can see is just K times the inverse covariance of the Gaussian density

Cramér-Rao Lower Bound

- the CRLB then says

$$\text{cov}(\hat{\boldsymbol{\mu}}|\mathbf{x}_{\text{meas}}) \geq \frac{1}{K}\boldsymbol{\Sigma} \quad (60)$$

- in other words, the lower limit of the uncertainty in the estimate of the mean, $\hat{\boldsymbol{\mu}}$, becomes smaller and smaller the more measurements we have (as we would expect)
- note, in computing the CRLB we did not need to actually specify the form of the unbiased estimator at all; **the CRLB is the lower bound for any unbiased estimator**

Example unbiased estimator: mean...

- in this case, it is not hard to find an estimator that performs right at the CRLB:

$$\hat{\boldsymbol{\mu}} = \frac{1}{K} \sum_{k=1}^K \mathbf{x}_{\text{meas},k} \quad (61)$$

- for the mean of this estimator we have

$$E[\hat{\boldsymbol{\mu}}] = E\left[\frac{1}{K} \sum_{k=1}^K \mathbf{x}_k\right] = \frac{1}{K} \sum_{k=1}^K E[\mathbf{x}_k] = \frac{1}{K} \sum_{k=1}^K \boldsymbol{\mu} = \boldsymbol{\mu} \quad (62)$$

which shows the estimator is indeed unbiased

...and the covariance

- for the covariance we have

$$\text{cov}(\hat{\boldsymbol{\mu}}|\mathbf{x}_{\text{meas}}) \quad (63a)$$

$$= E [(\hat{\boldsymbol{\mu}} - \boldsymbol{\mu})(\hat{\boldsymbol{\mu}} - \boldsymbol{\mu})^T] \quad (63b)$$

$$= E \left[\left(\frac{1}{K} \sum_{k=1}^K \mathbf{x}_k - \boldsymbol{\mu} \right) \left(\frac{1}{K} \sum_{k=1}^K \mathbf{x}_k - \boldsymbol{\mu} \right)^T \right] \quad (63c)$$

$$= \frac{1}{K^2} \sum_{k=1}^K \sum_{\ell=1}^K \underbrace{E [(\mathbf{x}_k - \boldsymbol{\mu})(\mathbf{x}_\ell - \boldsymbol{\mu})^T]}_{\Sigma \text{ when } k = \ell, \mathbf{0} \text{ otherwise}} \quad (63d)$$

$$= \frac{1}{K} \Sigma \quad (63e)$$

which is right at the CRLB; we can do no better

References

- Bayes, T., "Essay towards solving a problem in the doctrine of chances,"
Philosophical Transactions of the Royal Society of London, 1764.
- Thrun, S., Burgard, W., and Fox, D., *Probabilistic Robotics*, MIT Press, 2006.