Reasoning and Decision Making under Uncertainty

Classification with Context and Sensor Fusion

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Content and Overview

- Classification with Context
 - Viterbi Algorithm
- Hidden Markov Models
 - Concepts: Filtering, Prediction, Smoothing, Decoding, Learning
- Recursive Density Estimation
 - Kalman Filter
 - Particle Filter

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- Basic idea of "context"
 - Intuitively, it is obvious that classification in the context of multiple patterns should be possible with smaller error than classifying each pattern individually
 - Speech signals
 - Handwritten text
 - Other patterns embedded in a temporal or spatial context
 - We now consider a sequence of features

$$C = ({}^{1}c, {}^{2}c, \dots, {}^{N}c)$$

and the corresponding sequence of classes

$$\Omega = ({}^{1}\Omega, {}^{2}\Omega, \dots, {}^{N}\Omega) \text{ with } {}^{i}\Omega \in \{\Omega_{1}, \Omega_{2}, \dots, \Omega_{k}\}$$

ullet We are looking for the best sequence of classes taking into account all N decisions made

Obviously the Bayes classifier can be applied to a sequence of features and classes

$$p(\mathbf{\Omega}|\mathbf{C}) = \frac{p(\mathbf{\Omega})p(\mathbf{C}|\mathbf{\Omega})}{p(\mathbf{C})}$$

- Problem here: computational complexity
 - There are k^N values of the a priori probabilities and k^N densities of nN-dimensional vectors
 - First simplification: Feature vectors are statistically independent

$$p(\boldsymbol{C}|\boldsymbol{\Omega}) = \prod_{\rho=1}^{N} p(^{\rho}\boldsymbol{c}|^{\rho}\boldsymbol{\Omega} = \Omega_{\kappa})$$

Reduces to k densities of dimension n instead of k^N densities of nN-dimensional vectors

Second simplification: Class depends only on the direct predecessor (Markov property)

$$p(\mathbf{\Omega}) = p(^{1}\Omega, ^{2}\Omega, \dots, ^{N}\Omega) = p(^{1}\Omega)p(^{2}\Omega|^{1}\Omega)p(^{3}\Omega|^{1}\Omega^{2}\Omega) \dots p(^{N}\Omega|^{1}\Omega \dots ^{N-1}\Omega)$$
$$p(\mathbf{\Omega}) = p(^{1}\Omega)p(^{2}\Omega|^{1}\Omega)p(^{3}\Omega|^{2}\Omega) \dots p(^{N}\Omega|^{N-1}\Omega)$$

• Only k^2 transition probabilities and k probabilities (instead of k^N) left

Finally, the initial equation "simplifies" to

$$p(\mathbf{\Omega}|\mathbf{C}) = \frac{1}{p(\mathbf{C})}p(\mathbf{\Omega})p(\mathbf{C}|\mathbf{\Omega}) = \eta p(^{1}\Omega) \left(\prod_{\rho=2}^{N} p(^{\rho}\Omega|^{\rho-1}\Omega) \right) \left(\prod_{\rho=1}^{N} p(^{\rho}\mathbf{c}|^{\rho}\Omega = \Omega_{\kappa}) \right)$$

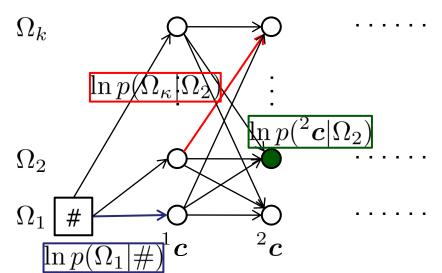
Better numerical properties are obtained using the logarithmized version

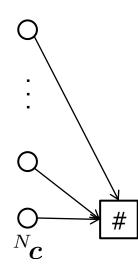
$$\ln p(\mathbf{\Omega}|\mathbf{C}) = \ln p(^{1}\Omega) + \sum_{\rho=2}^{N} \ln p(^{\rho}\Omega|^{\rho-1}\Omega) + \sum_{\rho=1}^{N} \ln p(^{\rho}\mathbf{c}|^{\rho}\Omega = \Omega_{\kappa}))$$

Do we know these densities and their parameters?

$$p({}^{\rho}\boldsymbol{c}|\Omega_{\kappa}) \qquad p(\Omega_{\kappa}|\Omega_{\lambda})$$

- Solution of the context problem: Viterbi algorithm
 - Set up network for all feature vectors and all possible classes: kN nodes
 - Each node represents observation o a feature vector under the assumption of a class
 - Assign weights to all nodes: $\ln p({}^{\rho}\boldsymbol{c}|\Omega_{\kappa})$
 - Each edge between two nodes represents the assumption that one class follows another class
 - Assign weights to all edges: $\ln p(\Omega_{\kappa}|\Omega_{\lambda})$
 - Sequence of features is delimited by a special symbol #
 - Can assign weights to start and end of sequence
 - lacktriangle Each sequence of classes for the sequence C corresponds to a path through the net
 - Weight of the path is sum of edge and node weights
 - Optimal path through the net is the one with maximum weight
 - Viterbi algorithm calculates the optimal path using principle of dynamic programming





Viterbi algorithm [Vit67]

FOR $\kappa = 1$ TO NFOR $\lambda = 1$ TO k

Calculate
$$G_{\lambda,\rho} = \ln p({}^{\rho}\boldsymbol{c}|\Omega_{\lambda}) + \max_{j \in \{1,...,k\}} (G_{j,\rho-1} + \ln p(\Omega_{\lambda}|\Omega_{j}))$$

Special treatment for $\rho = 1$

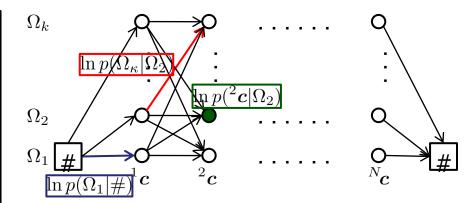
$$\max_{j} (G_{j,\rho-1} + \ln p(\Omega_{\lambda}|\Omega_{j})) = \ln p(\Omega_{\lambda}|\#) = \ln p(\Omega_{\lambda})$$

Store index of the predecessor node leading to the maximum of $G_{\lambda,\rho}$

The last column contains k weights. Calculate largest weight

$$G_{\kappa,N} = \max_{\lambda} G_{\lambda,N}$$

Determine path with maximum weight by tracing stored indexes from end node to start node: Sequence of these nodes is $\,\Omega\,$



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What we came across with the Viterbi algorithm is better known under the term

(Hidden) Markov Model (HMM)

- Getting closer to the common view of HMMs we need to broaden the mathematical perspective
- Change of notation from now on
 - We generalize our features to observations $c \rightarrow o$
 - We generalize classes to states (may be continuous) $\Omega \rightarrow q$
 - States can only be estimated, they can never be observed, they are hidden (latent variables)
 - States differ in some important properties from classes
 - A state is usually multivariate and sometimes continuous
 - A state represents a more comprehensive description of the world
- Notation of time-sequential states and observations

$$\langle \boldsymbol{q} \rangle_t = \boldsymbol{q}_t, \boldsymbol{q}_{t-1}, \dots, \boldsymbol{q}_0$$

$$\langle \boldsymbol{o} \rangle_t = \boldsymbol{o}_t, \boldsymbol{o}_{t-1}, \dots, \boldsymbol{o}_0$$

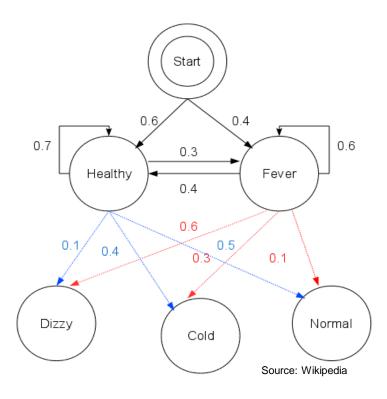
The context equation now changes to

$$p(\boldsymbol{q}_{t}, \boldsymbol{q}_{t-1}, \dots, \boldsymbol{q}_{0} | \boldsymbol{o}_{t}, \boldsymbol{o}_{t-1}, \dots, \boldsymbol{o}_{0}) = p(\langle \boldsymbol{q}_{t} \rangle | \langle \boldsymbol{o}_{t} \rangle)$$

$$= \frac{p(\langle \boldsymbol{q} \rangle_{t}) p(\langle \boldsymbol{o} \rangle_{t} | \langle \boldsymbol{q} \rangle_{t})}{p(\langle \boldsymbol{o} \rangle_{t})}$$

$$= \eta \prod_{t} p(\boldsymbol{q}_{t} | \boldsymbol{q}_{t-1}) \prod_{t} p(\boldsymbol{o}_{t} | \boldsymbol{q}_{t})$$

- lacksquare What about $p(oldsymbol{q}_0|oldsymbol{q}_{-1})$?
- In the context of HMM the "transition probabilities" are based mostly on discrete states
 - Just like our discrete classes from above
 - This leads to a transition matrix that contains all the transition probabilities
 - But: we do not want to restrict to discrete states
- In the context of HMM the observations are discrete most of the time
 - At each time any of the (finite) observations can occur
 - This leads to a observation matrix that contains the probabilities for a observation under the condition of a specific current state
 - But: we do not want to restrict to discrete observations



- Known states: s_1 = Healthy, s_2 = Fever
- Known observations: v_1 = Dizzy, v_2 = Cold, v_3 = Normal
- Start probabilities

$$\pi = (0.6 \quad 0.4) \qquad \pi_i = p(q_0 = s_i)$$

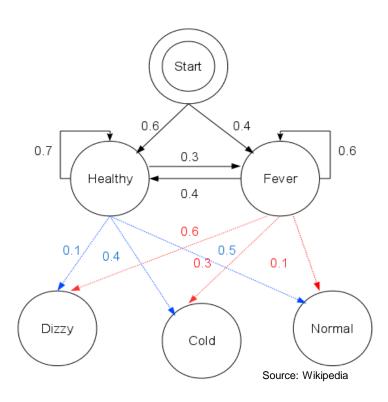
State transition probabilities written as matrix

$$T = \begin{pmatrix} 0.7 & 0.3 \\ 0.6 & 0.4 \end{pmatrix} \qquad t_{ij} = p(q_t = s_j | q_{t-1} = s_i)$$

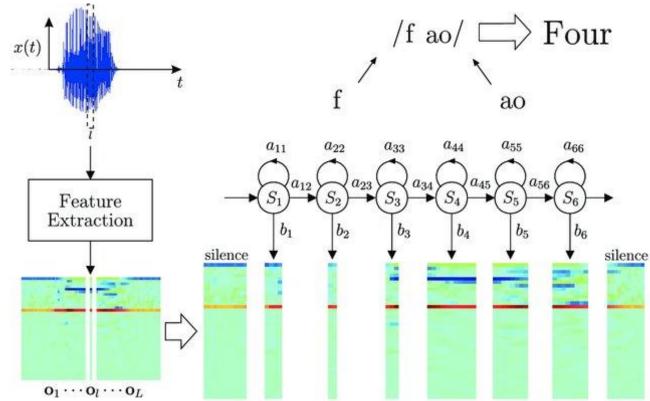
Output probabilities (emission probabilities) written as matrix

$$\mathbf{E} = \begin{pmatrix} 0.1 & 0.4 & 0.5 \\ 0.6 & 0.3 & 0.1 \end{pmatrix} \qquad e_{ij} = p(o_t = v_j | q_t = s_i) = e_i(j)$$

- **Example state sequence** $\langle q \rangle_4 = ($ Healthy, Healthy, Fever, Fever, Healthy)
- Example observation sequence $\langle o \rangle_4 = (\text{Cold, Dizzy, Dizzy, Normal, Cold})$



- Classification with discrete observations and discrete states makes sense for many applications
 - Example: speech recognition
 - Observation in HMM model: short-term spectra of the speech signal
 - (Hidden) state in HMM model: semantic units (e.g. phonemes in speech recognition)
 - HMM model used to detect sequence of semantic units in the sequential data



Important HMM concepts

- Filtering problem
 - ullet Given a HMM and an observation sequence of length T
 - Looking for the probability that the instantaneous hidden state at the last time T is some specific state
 - An efficient method for solving the filtering problem is the forward algorithm
- Prediction problem
 - ullet Given a HMM and an observation sequence of length T
 - We are looking for the probability that the HMM will be in a specific hidden state at future time T+t (t time-steps ahead in the future)
 - Prediction is repeated filtering without new observations
 - An efficient method for solving the prediction problem is the forward algorithm
- Smoothing problem
 - ullet Given a HMM and an observation sequence of length $\,T\,$
 - ullet We are looking for the probability that the model was in a certain state at a previous time T t
 - Smoothing uses future observations
 - An efficient method for solving the smoothing problem is the forward-backward algorithm

Decoding problem

- Given an HMM and an observation sequence
- Determine the most probable sequence of states that could have produced a given output sequence
- An efficient method for solving the decoding problem is the Viterbi algorithm
- Learning problem
 - Given only an observation sequence
 - Determine the parameters of an HMM that are most likely to produce the output sequence
 - An efficient method for solving the decoding problem is the Baum-Welch algorithm (specialized EM-Algorithm)

- Forward algorithm
 - Probability in the given HMM to make series of observations $\langle o \rangle_T$
 - 1. Initialization for all known states j at first time step 0

$$\alpha_0(j) = \pi_j e_j(o_0)$$

2. Recursion for a all known states j and time steps t

$$\alpha_t(j) = e_j(o_t) \sum_i \alpha_{t-1}(i) t_{ij}$$

3. Termination

$$p(\langle o \rangle_T | \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E}) = \sum_i \alpha_T(i)$$

$$\pi_i = p(q_0 = s_i)$$
 $t_{ij} = p(q_t = s_j | q_{t-1} = s_i)$
 $e_{ij} = p(o_t = v_j | q_t = s_i) = e_i(j)$

- Backward algorithm
 - lacktriangle Probability of partial observation sequence from t+1 to end T, given state at time t
 - 1. Initialization for all known states i at last time step T

$$\beta_T(i) = 1 \quad \forall i$$

2. Recursion for a all known states i and time steps t

$$\beta_t(i) = \sum_j t_{ij} e_j(o_{t+1}) \beta_{t+1}(j)$$

3. Termination

$$p(\langle o \rangle_T | \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E}) = \sum_j \pi_j e_j(o_0) \beta_0(j)$$

$$\pi_i = p(q_0 = s_i)$$
 $t_{ij} = p(q_t = s_j | q_{t-1} = s_i)$
 $e_{ij} = p(o_t = v_j | q_t = s_i) = e_i(j)$

- Using forward and backward variables for solving the smoothing problem
 - Probability that the model was in a certain state at a previous time if complete sequence of observations is available
 - Use combination of forward and backward variables
 - lacktriangle Matrix S aggregates these probabilities for all states and time steps

$$s_{it} = \alpha_t(i)\beta_t(i) = s_t(i)$$

- Viterbi algorithm
 - Find the most probable sequence of states given a series of observations
 - 1. Initialization for all known states j at first time step 0

$$v_0(j) = \pi_j e_j(o_0)$$
$$b_0(j) = 0$$

2. Recursion for a all known states j and time steps t

$$v_t(j) = \max_{i} v_{t-1}(i)t_{ij}e_j(o_t)$$
$$b_t(j) = \operatorname*{argmax}_{i} v_{t-1}(i)t_{ij}e_j(o_t)$$

- Termination
 - Best possible score: $\max_{i} v_T(i)$
 - Start of backtrace: $\underset{i}{\operatorname{argmax}} v_T(i)$

$$\pi_i = p(q_0 = s_i)$$
 $t_{ij} = p(q_t = s_j | q_{t-1} = s_i)$
 $e_{ij} = p(o_t = v_j | q_t = s_i) = e_i(j)$

- Baum-Welch Algorithm
 - Finds/calculates the parameters π , T, E of the HMM
 - Initialization parameters π , T, E of the HMM randomly or with prior knowledge
 - Compute forward and backward variables: $\alpha_t(j), \beta_t(i)$
 - Remember that

$$p(\langle o \rangle_T | \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E}) = \sum_i \alpha_T(i) = \sum_j \pi_j e_j(o_0) \beta_0(j)$$

$$\pi_i = p(q_0 = s_i)$$
 $t_{ij} = p(q_t = s_j | q_{t-1} = s_i)$
 $e_{ij} = p(o_t = v_j | q_t = s_i) = e_i(j)$

- EM-Algorithm: Iterate until convergence
 - E-step
 - lacktriangle Probability of being in state s_i at time t, given observation sequence and model

$$\gamma_t(i) = p(q_t = s_i | \langle o \rangle_T, \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E}) = \frac{\alpha_t(i)\beta_t(i)}{p(\langle o \rangle_T | \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E})}$$

$$\pi_i = p(q_0 = s_i)$$
 $t_{ij} = p(q_t = s_j | q_{t-1} = s_i)$
 $e_{ij} = p(o_t = v_j | q_t = s_i) = e_i(j)$

• Probability of being in state s_i at time t and state s_j at time t+1, given observation sequence and model

$$\xi_t(i,j) = p(q_t = s_i, q_{t+1} = s_j | \langle o \rangle_T, \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E}) = \frac{\alpha_t(i) t_{ij} e_j(o_{t+1}) \beta_{t+1}(j)}{p(\langle o \rangle_T | \boldsymbol{\pi}, \boldsymbol{T}, \boldsymbol{E})}$$

This satisfies the relationship

$$\gamma_t(i) = \sum_j \xi_t(i,j)$$

- M-step
 - Update state transition probabilities

$$t_{ij} = \frac{\sum_{t=0}^{T-1} \xi_t(i,j)}{\sum_{t=0}^{T-1} \sum_{k} \xi_t(i,k)}$$

Update emission probabilities

$$e_j(k) = \frac{\sum_{t=0}^{T} f(o_t, v_k) \gamma_t(j)}{\sum_{t=0}^{T} \gamma_t(j)}$$

$$f(a, b) = \begin{cases} 1 & \text{if } a = b \\ 0 & \text{otherwise} \end{cases}$$

Convergence criteria: maximize likelihood

- Baum-Welch algorithm so far estimates parameters from only one sequence of observations
 - Many observation sequences are used for practical implementations
 - E-step: Estimate γ and ξ for every single observation
 - M-step: Average over updates for transition and emission probabilities
- All terms in the estimation process are significantly less than 1.0
 - Products and summations converge to 0.0 very quickly (the larger T becomes, the faster the calculations will converge towards 0.0)
 - Precision of computation of forward/backward variables will exceed the floating-point accuracy of CPU
 - Solution is a proper scaling of forward and backward variables
 - See eLearning: Dawei Shen. "Some Mathematics for HMM"

Classification with Context – Sensor Fusion

- Disadvantages Forward Algorithm and Viterbi Algorithm
 - All observations must be available at the time of calculation
 - Disadvantage 1: Stepwise integration of new information not possible

$$p(\langle \boldsymbol{q}_t \rangle | \langle \boldsymbol{o}_t \rangle) = \eta \prod_t p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) \prod_t p(\boldsymbol{o}_t | \boldsymbol{q}_t)$$

- Need to have all observations available
- Calculation goes over complete sequence of observations
- Disadvantage 2: Need to find proper family of densities that describes scenario
 - In common HMM applications this is often achieved by discrete transitions and discrete observations

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Recursive Density Estimation

- More general approach: recursive density estimation
 - Start again with same Bayes classifier as before
 - Limit our interest to an estimate of the state at the current time

$$p(\langle \boldsymbol{q} \rangle_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t)$$

- Previous states have already been calculated
- This can be rewritten as

$$p(\langle \boldsymbol{q} \rangle_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$$

Recursive Density Estimation

Components of the recursive formulation



- Estimate density for state at current time: Old information is not of interest in dynamic environment
- Recursive part contains all information up to the last time step. Recursion until time step 0 is reached
- State transition contains information about system dynamics
- Observation integrates new information/observations/patterns
- Disadvantage 1 solved: Stepwise integration of new information now possible
 - Only density from last time step, system dynamics (known) and current observation is needed at any time

Recursive Density Estimation

Disadvantage 2: There are different solutions

$$p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$$

- It is always only a matter of implementing the above formula
- We are looking for ways to describe these densities
- Will discuss two techniques
 - Kalman Filter
 - Particle Filter

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- ▶ Realization of Recursive Density Estimation: Kalman Filter
 - Dynamic environments rely on interaction with the environment
 - lacktriangle Introduce (control) actions a in addition to observations
 - Extend recursive formula by actions

$$p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t, \langle \boldsymbol{a} \rangle_{t-1}) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}, \boldsymbol{a}_{t-1}) p(\boldsymbol{q}_{t-1} | \langle \boldsymbol{o} \rangle_{t-1}, \langle \boldsymbol{a} \rangle_{t-2}) d\boldsymbol{q}_{t-1}$$

- State is thus dependent on
 - Observations: passive component, has no own influence on the environment, only observes the environment
 - Actions: our own commands for interaction with the environment, own decision, if no actions exist then omit them
- Kalman filter has two major requirements
 - Linear state transition
 - Normally distributed densities

$$p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t, \langle \boldsymbol{a} \rangle_{t-1}) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}, \boldsymbol{a}_{t-1}) p(\boldsymbol{q}_{t-1} | \langle \boldsymbol{o} \rangle_{t-1}, \langle \boldsymbol{a} \rangle_{t-2}) d\boldsymbol{q}_{t-1}$$

- Kalman requirement: linear state transition
 - Realization of

$$p(\boldsymbol{q}_t|\boldsymbol{q}_{t-1},\boldsymbol{a}_{t-1})$$

with linear state transition

$$oldsymbol{q}_t = oldsymbol{A}_t oldsymbol{q}_{t-1} + oldsymbol{B}_t oldsymbol{a}_{t-1} + oldsymbol{\epsilon}_t$$

Normally distributed random variable with covariance matrix Q

and normally distributed uncertainty leads to density:

$$p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}, \boldsymbol{a}_{t-1}) = \frac{1}{\sqrt{\det(2\pi\boldsymbol{Q}_t)}} \exp\left(-\frac{(\boldsymbol{q}_t - (\boldsymbol{A}_t \boldsymbol{q}_{t-1} + \boldsymbol{B}_t \boldsymbol{a}_{t-1}))^T \boldsymbol{Q}_t^{-1} (\boldsymbol{q}_t - (\boldsymbol{A}_t \boldsymbol{q}_{t-1} + \boldsymbol{B}_t \boldsymbol{a}_{t-1}))}{2}\right)$$

$$p(\boldsymbol{q}_{t}|\langle\boldsymbol{o}\rangle_{t},\langle\boldsymbol{a}\rangle_{t-1}) = p(\boldsymbol{o}_{t}|\boldsymbol{q}_{t}) \int p(\boldsymbol{q}_{t}|\boldsymbol{q}_{t-1},\boldsymbol{a}_{t-1}) p(\boldsymbol{q}_{t-1}|\langle\boldsymbol{o}\rangle_{t-1},\langle\boldsymbol{a}\rangle_{t-2}) d\boldsymbol{q}_{t-1}$$

- Kalman requirement: linear observation model
 - Realization of

$$p(\boldsymbol{o}_t|\boldsymbol{q}_t)$$

with linear observation model

$$oldsymbol{o}_t = oldsymbol{C}_t oldsymbol{q}_t + oldsymbol{\delta}_t$$

Normally distributed random variable with covariance matrix *R*

and normally distributed uncertainty leads to density:

$$p(\boldsymbol{o}_t|\boldsymbol{q}_t) = \frac{1}{\sqrt{\det(2\pi\boldsymbol{R}_t)}} \exp\left(-\frac{(\boldsymbol{o}_t - \boldsymbol{C}_t\boldsymbol{q}_t)^T \boldsymbol{R}_t^{-1} (\boldsymbol{o}_t - \boldsymbol{C}_t\boldsymbol{q}_t)}{2}\right)$$

- Famous Kalman algorithm [Kal60] is a realization of this linear state transition and linear observation model
 - Need mean vector and covariance matrix of a normal distribution representing

$$p(\boldsymbol{q}_t|\boldsymbol{q}_{t-1},\boldsymbol{a}_{t-1})$$

- Inputs for one iteration of the Kalman filter
 - Mean vector and covariance matrix of last iteration (last time step)
 - Current observation and action
- State at time 0 must be normally distributed
 - Need initial mean vector and covariance matrix

 $\mathtt{Kalman}(oldsymbol{\mu}_{t-1}, oldsymbol{\Sigma}_{t-1}, oldsymbol{o}_t, oldsymbol{a}_{t-1})$

1.
$$\mu'_t = A_t \mu_{t-1} + B_t a_{t-1}$$

2.
$$\boldsymbol{\Sigma}_t' = \boldsymbol{A}_t \boldsymbol{\Sigma}_{t-1} \boldsymbol{A}_t^T + \boldsymbol{Q}_t$$

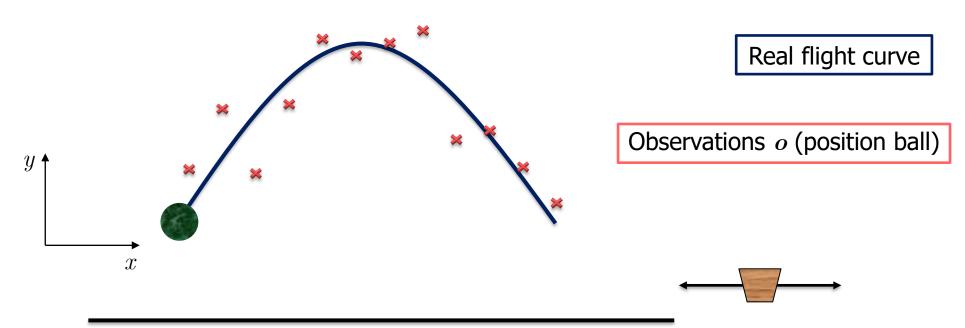
3.
$$\boldsymbol{K}_{t} = \boldsymbol{\Sigma}_{t}^{\prime} \boldsymbol{C}_{t}^{T} (\boldsymbol{C}_{t} \boldsymbol{\Sigma}_{t}^{\prime} \boldsymbol{C}_{t}^{T} + \boldsymbol{R}_{t})^{-1}$$
4. $\boldsymbol{\mu}_{t} = \boldsymbol{\mu}_{t}^{\prime} + \boldsymbol{K}_{t} (\boldsymbol{o}_{t} - \boldsymbol{C}_{t} \boldsymbol{\mu}_{t}^{\prime})$
5. $\boldsymbol{\Sigma}_{t} = (\boldsymbol{I} - \boldsymbol{K}_{t} \boldsymbol{C}_{t}) \boldsymbol{\Sigma}_{t}^{\prime}$

4.
$$\mu_t = \mu'_t + K_t(o_t - C_t \mu'_t)$$

5.
$$\Sigma_t = (\boldsymbol{I} - \boldsymbol{K}_t \boldsymbol{C}_t) \Sigma_t'$$

- Important component: Kalman gain K [Thr05]
 - Degree/Weighting with which observation is integrated
 - Minimizes the expected, a posteriori error
 - Solves question: Greater confidence in prediction or in measurement of observation?

Example Kalman filter: parabolic throw



Physical movement

$$x_{t+1} = x_t + v_x \Delta t$$
$$y_{t+1} = y_t + v_y \Delta t - 0.5g \Delta t^2$$

Required: transition and observation model

Transition model:

$$oldsymbol{q}_t = oldsymbol{A}_t oldsymbol{q}_{t-1} + oldsymbol{B}_t oldsymbol{a}_{t-1} + oldsymbol{\epsilon}_t$$

$$\boldsymbol{q}_{t} = \begin{pmatrix} x \\ y \\ v_{x} \\ v_{y} \end{pmatrix}_{t} = \begin{pmatrix} 1 & 0 & \Delta t & 0 \\ 0 & 1 & 0 & \Delta t \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix} \begin{pmatrix} x \\ y \\ v_{x} \\ v_{y} \end{pmatrix}_{t-1} + \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0.5\Delta t^{2} & 0 \\ 0 & 0 & 0 \\ 0 & \Delta t & 0 \end{pmatrix} \begin{pmatrix} 0 \\ -g \\ 0 \end{pmatrix}_{t-1} + \boldsymbol{\epsilon}_{t-1}$$

Observation model:

$$oldsymbol{o}_t = oldsymbol{C}_t oldsymbol{q}_t + oldsymbol{\delta}_t$$

$$oldsymbol{o}_t = \begin{pmatrix} x \\ y \end{pmatrix}_t = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \end{pmatrix} \begin{pmatrix} x \\ y \\ v_x \\ v_y \end{pmatrix}_t + oldsymbol{\delta}_t$$

Estimated mean vector corresponds to our state

$$\mu=q$$

Arbitrary initial mean vector, e.g.

$$\boldsymbol{\mu}_0 = (0, 0, 0, 0)^T$$

Arbitrary initial covariance matrix, e.g.

$$oldsymbol{\Sigma}_0 = egin{pmatrix} 1 & 0 & 0 & 0 \ 0 & 1 & 0 & 0 \ 0 & 0 & 1 & 0 \ 0 & 0 & 0 & 1 \end{pmatrix}$$

Action constant: acceleration due to gravity

$$\boldsymbol{a} = (0, -g, 0)^T$$

4. $\mu_t = \mu'_t + K_t(o_t - C_t \mu'_t)$

3. $\boldsymbol{K}_t = \boldsymbol{\Sigma}_t' \boldsymbol{C}_t^T (\boldsymbol{C}_t \boldsymbol{\Sigma}_t' \boldsymbol{C}_t^T + \boldsymbol{R}_t)^{-1}$

 $\mathtt{Kalman}(oldsymbol{\mu}_{t-1}, oldsymbol{\Sigma}_{t-1}, oldsymbol{o}_t, oldsymbol{a}_{t-1})$

1. $\mu'_t = A_t \mu_{t-1} + B_t a_{t-1}$

2. $\Sigma_t' = A_t \Sigma_{t-1} A_t^T + Q_t$

5. $\Sigma_t = (I - K_t C_t) \Sigma_t'$

lacksquare Information about noise in \emph{Q} and \emph{R} depend on type of observation

- Limitations of the Kalman filter
 - Linear state transition and observation: unrealistic/wrong in most applications
 - Example: Robot moving on circular path cannot be described with linear motion model
 - Only unimodal, Gaussian distributed states can be modeled
 - Noise component limited

- Extended Kalman Filter (EKF)
 - lacktriangle Eliminates linearity requirement: nonlinear functions g and h

Kalman EKF
$$m{q}_t = m{A}_tm{q}_{t-1} + m{B}_tm{a}_{t-1} + m{\epsilon}_t \ m{o}_t = m{C}_tm{q}_t + m{\delta}_t \ m{o}_t = m{h}(m{q}_t) + m{\delta}_t$$

- EKF idea: Linearization
 - Calculate a function that locally approximates a nonlinear function
 - There are many methods for linearization
 - Common idea: Take first both terms of the Taylor series expansion
 - lacksquare Leads to usage of Jacobian G and H of the functions g and h

$$\mathtt{Kalman}(oldsymbol{\mu}_{t-1}, oldsymbol{\Sigma}_{t-1}, oldsymbol{o}_t, oldsymbol{a}_{t-1})$$

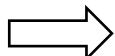
1.
$$\mu'_t = A_t \mu_{t-1} + B_t a_{t-1}$$

2.
$$\Sigma_t' = A_t \Sigma_{t-1} A_t^T + Q_t$$

3.
$$\boldsymbol{K}_t = \boldsymbol{\Sigma}_t' \boldsymbol{C}_t^T (\boldsymbol{C}_t \boldsymbol{\Sigma}_t' \boldsymbol{C}_t^T + \boldsymbol{R}_t)^{-1}$$

4.
$$\mu_t = \mu'_t + K_t(o_t - C_t \mu'_t)$$

5.
$$\Sigma_t = (\boldsymbol{I} - \boldsymbol{K}_t \boldsymbol{C}_t) \Sigma_t'$$



EKF
$$(oldsymbol{\mu}_{t-1}, oldsymbol{\Sigma}_{t-1}, oldsymbol{o}_t, oldsymbol{a}_{t-1})$$

$$1. \ oldsymbol{\mu}_t' = oldsymbol{g}(oldsymbol{\mu}_{t-1}, oldsymbol{a}_{t-1})$$

1.
$$\mu'_t = g(\mu_{t-1}, a_{t-1})$$

2.
$$\Sigma_t' = G_t \Sigma_{t-1} G_t^T + Q_t$$

3.
$$\boldsymbol{K}_{t} = \boldsymbol{\Sigma}_{t}^{\prime} \boldsymbol{H}_{t}^{T} (\boldsymbol{H}_{t} \boldsymbol{\Sigma}_{t}^{\prime} \boldsymbol{H}_{t}^{T} + \boldsymbol{R}_{t})^{-1}$$
4. $\boldsymbol{\mu}_{t} = \boldsymbol{\mu}_{t}^{\prime} + \boldsymbol{K}_{t} (\boldsymbol{o}_{t} - \boldsymbol{h}(\boldsymbol{\mu}_{t})^{\prime})$
5. $\boldsymbol{\Sigma}_{t} = (\boldsymbol{I} - \boldsymbol{K}_{t} \boldsymbol{H}_{t}) \boldsymbol{\Sigma}_{t}^{\prime}$

4.
$$\mu_t = \mu'_t + K_t(o_t - h(\mu_t)')$$

5.
$$\Sigma_t = (\boldsymbol{I} - \boldsymbol{K}_t \boldsymbol{H}_t) \Sigma_t'$$

- EKF still has serious limitations
 - Only unimodal, Gaussian distributed states can be modeled
 - Noise component limited

- There are many other methods that deal with these constraints
 - Unscented Kalman filter
 - For highly nonlinear state transitions and observation
 - Samples around mean vector
 - Information filter
 - Histogram filter
 - Particle filter

Content and Overview

- Classification with Context
 - Viterbi Algorithm
- Hidden Markov Models
 - Concepts: Filtering, Prediction, Smoothing, Decoding, Learning
- Recursive Density Estimation
 - Kalman Filter
 - Particle Filter

- Realization recursive density estimation: Particle Filter
 - Can approximate any kind of complex densities
 - Not restricted to normal distributions
 - Non-parametric
 - Suitable for nonlinear problems
 - Solution possibility for recursive density estimation
 - Need to look for solutions for two questions and find suitable components
 - Task 1: Unlimited possibility to model arbitrary densities
 - Solution: Particle Sets
 - Particle sets model densities
 - Task 2: Any system dynamics should be describable
 - Solution: Particle Filter
 - Particle filters model system dynamic

- Task 1: Approximate single density by particle set
 - Particle set consists of individual particles

$$S_t = \{s_t^1, s_t^2, \dots, s_t^n\}$$

• Each particle contains necessary data s for modeling the state q, e.g. 2d position

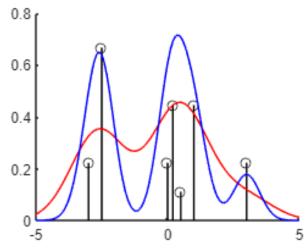
$$s = q = (x, y)^T$$
 $s = \{s, w\} = \{(x, y)^T, w\}$

- Each particle contains state information: application-dependent, usually same as q
- Each particle contains weighting w (think of it as probability of this particle)
- Each particle describes an evaluation of the density at a location in the state space
- Approximation of complete density from particle set

$$p(\mathbf{q}_t | \langle \mathbf{o} \rangle_t) \approx \sum_i w_t^i \delta\left(\mathbf{s}_t^i - \mathbf{q}_t\right)$$

$$\sum_i w_t^i = 1$$

- Very similar to Parzen estimate, only additional weighting for particle (like in mixture distributions)
- If number of particles runs to infinity, window function δ can be arbitrary small and density can be approximated with arbitrary accuracy



- Task 2: Particle filter that handles system dynamics over time
 - Most important particle filter: Condensation Algorithm [Isa98]

Condensation Algorithm

- 1. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from S_{t-1} for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)

There are many other particle filters [Dou01]

• Initial particle set S_0

Condensation Algorithm

- 1. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from $S_{t extstyle{-}1}$ for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)
- lacktriangle Mostly no prior knowledge about initial state available: Uniform distribution assumption for $p(m{q}_0)$
 - lacktriangle Create initial particle set by distributing n particles uniformly over state space
- If prior knowledge available: Draw n particles from this distribution
- Set particle weights $w_0^i = \frac{1}{n}$

• Sample n particles from S_{t-1} and create new particle set S_t

Condensation Algorithm

- I. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from $S_{t extstyle{-}1}$ for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)

How to sample new particles

- Create uniformly distributed random number [0.0 1.0]
- Find particles with matching cumulative range (random number within range)
- Copy particles into new particle set
- This is the same as sampling from multinominal distribution

Particle	Weight	Cumulative Range
s_1	0.1	0.0 – 0.1
\$2	0.3	0.1 – 0.4
S_3	0.45	0.4 – 0.85
s_4	0.15	0.85 – 1.0

Example:

- 1. Created random number: 0.92
- 2. Created random number: 0.58
- 3. Created random number: 0.89
- 4. Created random number: 0.27

$S_t = \{s_4, s_3, s_4, s_2\}$	

New particle set

- Propagate particles
 - State transition model

$$p(m{q}_t|m{q}_{t-1},m{a}_{t-1})$$
 or $p(m{q}_t|m{q}_{t-1})$ or $p(m{q}_t|m{q}_{t-1},m{o}_{t-1})$ or ...

- **Condensation Algorithm**
- 1. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from $S_{t ext{-}1}$ for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)

- Goal: Adapt state data in particle due to (known or assumed) system dynamics
- In general, state transition is not deterministic: sample from appropriate density

$$p(\langle \boldsymbol{q} \rangle_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$$

Evaluate particle

$$p(\boldsymbol{o}_t|\boldsymbol{q}_t)$$

- **Condensation Algorithm**
- 1. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from S_{t-1} for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)

- Set weight of each particle
 - Calculate new particle weight with classifier
 - Example for simple normal distribution classifier

$$w_t^i = p\left(\boldsymbol{o}_t | (x_t^i, y_t^i)^T\right) = \mathcal{N}\left(\boldsymbol{o}_t | (x_t^i, y_t^i)^T, \boldsymbol{\Sigma}\right)$$

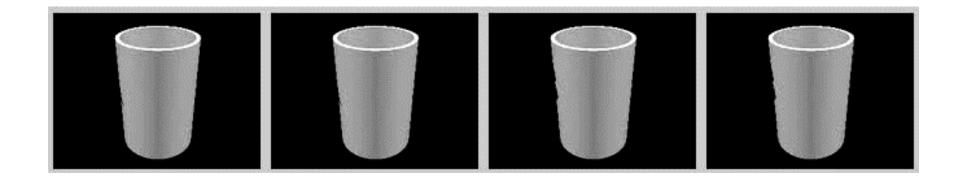
 $S_t = \{s_t^1, s_t^2, \dots, s_t^n\}$ $s_t^i = \{(x_t^i, y_t^i)^T, w_t^i\}$

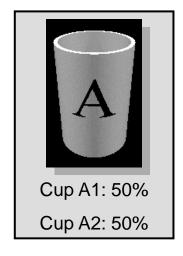
- Usually: Use your statistical classifier here
 - Calculate probability for each sample for the current, given observation
 - This is like evaluating density for feature vector o
- After evaluating all particles: Normalize weighting

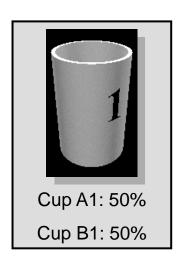
$$w_t^i = \frac{w_t^i}{\sum_j w_t^j}$$

$p(\langle \boldsymbol{q} \rangle_t \, | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \, \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$ Particle Filter — Classification/Localization Example

- Example particle filter: Classification
 - Recognize ambiguous objects [Dei01]

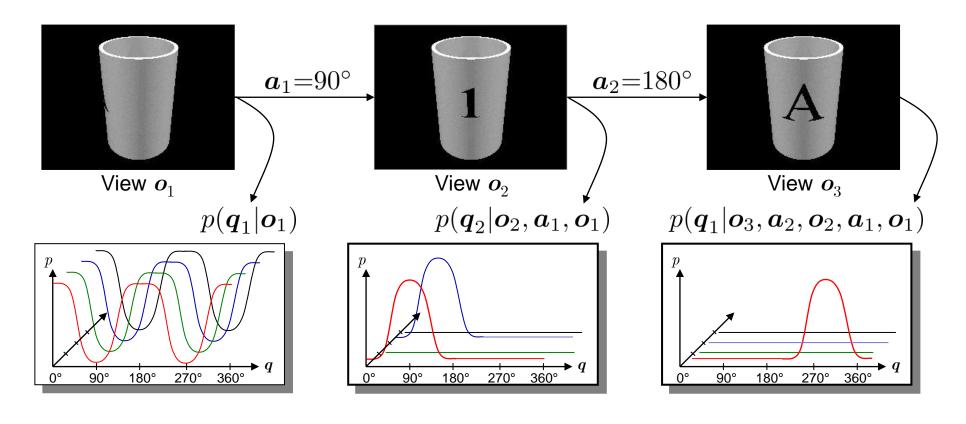






$p(\langle \boldsymbol{q} \rangle_t \, | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \, \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$ Particle Filter — Classification/Localization Example

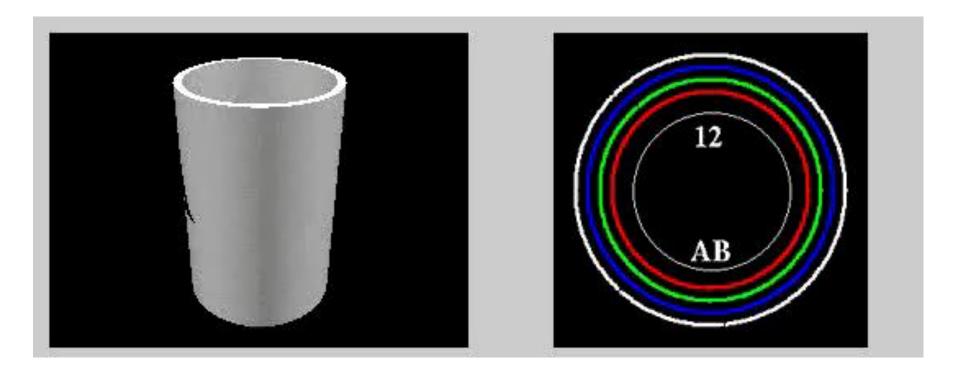
- Example particle filter: Classification
 - Fusion of information from 3 images



Cup A1, Cup A2, Cup B1, Cup B2

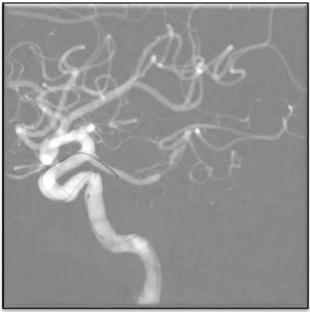
$$p(\langle \boldsymbol{q} \rangle_t \, | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \, \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$$
Particle Filter — Classification/Localization Example

- Example particle filter: Classification
 - Fusion of a sequence of many images [Dei01]



- Example particle filter: Medical Application in Angiography
 - 2-D X-ray imaging basis of many applications in interventional radiology
 - Specialized C-arm systems used for treatment of strokes, heart attacks, tumors

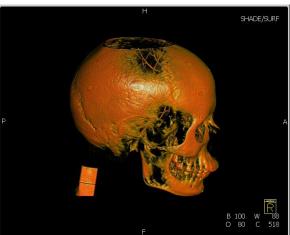






- Example particle filter: Medical Application in Angiography
 - 3-D interventional imaging essential technique nowadays
 - Information from 3D volume often not visible in 2D X-ray images
 - No real-time 3D imaging





3D volume

- Example particle filter: Medical Application in Angiography
 - Application: Overlay of rendered 3D volume
 - Preprocessing of 3D volume (e.g. remove
 - unwanted anatomic details) possible
 - Overlaid volume will follow any C-arm system parameter changes and movements
 - Here: Treatment of an aneurysm



$p(\langle \boldsymbol{q} \rangle_t \,|\, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t |\, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} |\, \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$

Particle Filter – Medical Example (Angiography)

- Example particle filter: Medical Application in Angiography
 - Track and reconstruct medical guidewire in patient vessel system [Brü09]



Sequence of X-ray images

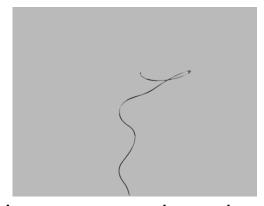
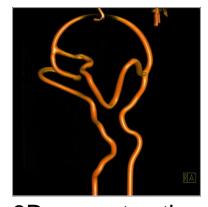
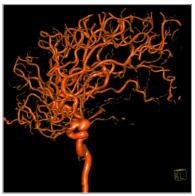


Image processing using mask images: only guidewire visible

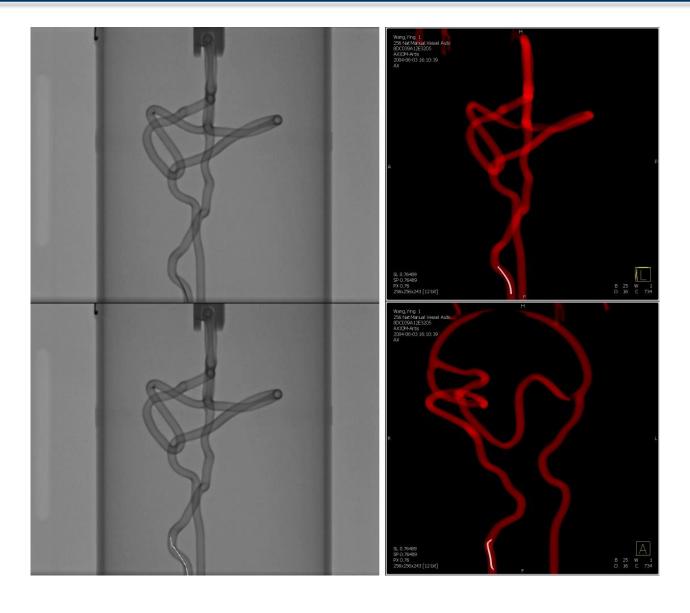


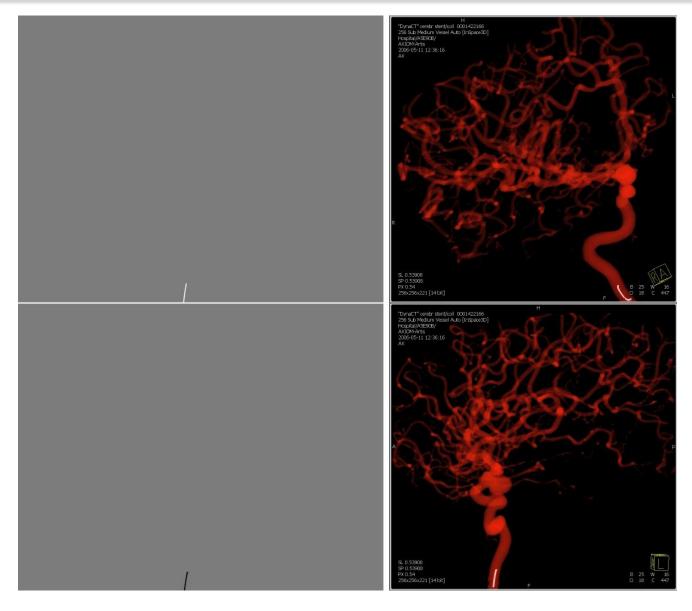
3D reconstruction from CT-scanner

 Difficult problem in case of complex vessel systems

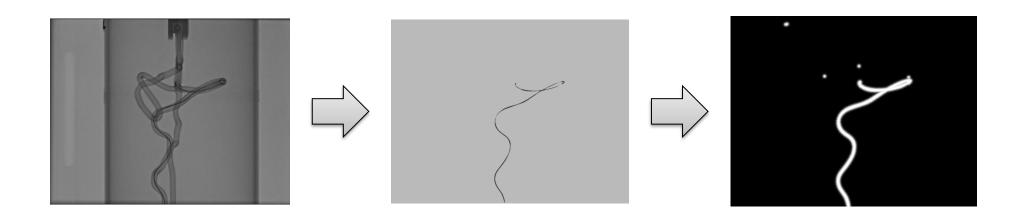


$\begin{array}{c} p(\langle \boldsymbol{q} \rangle_t \, | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \, \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1} \\ \textbf{Particle Filter - Medical Example (Angiography)} \end{array}$





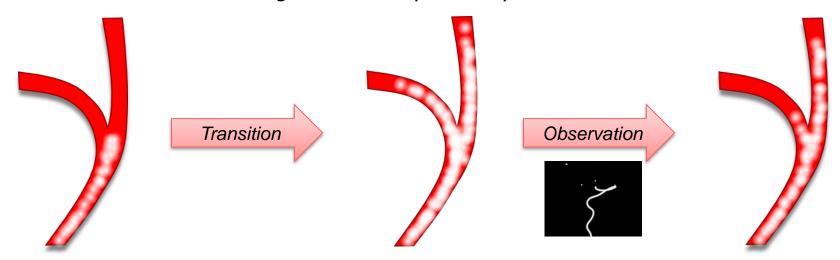
- Example particle filter: Catheter Tracking
 - Use 3D coordinates as state $q = (x, y, z)^T$
 - Observation $p(\boldsymbol{o}_t|\boldsymbol{q}_t)$
 - Use image difference (live image mask image) o_t
 - Apply some kind of blurring → model inaccuracies
 - Density $p(o_t|q_t)$ of observations depend on known projective geometry (just like for catheter tip)



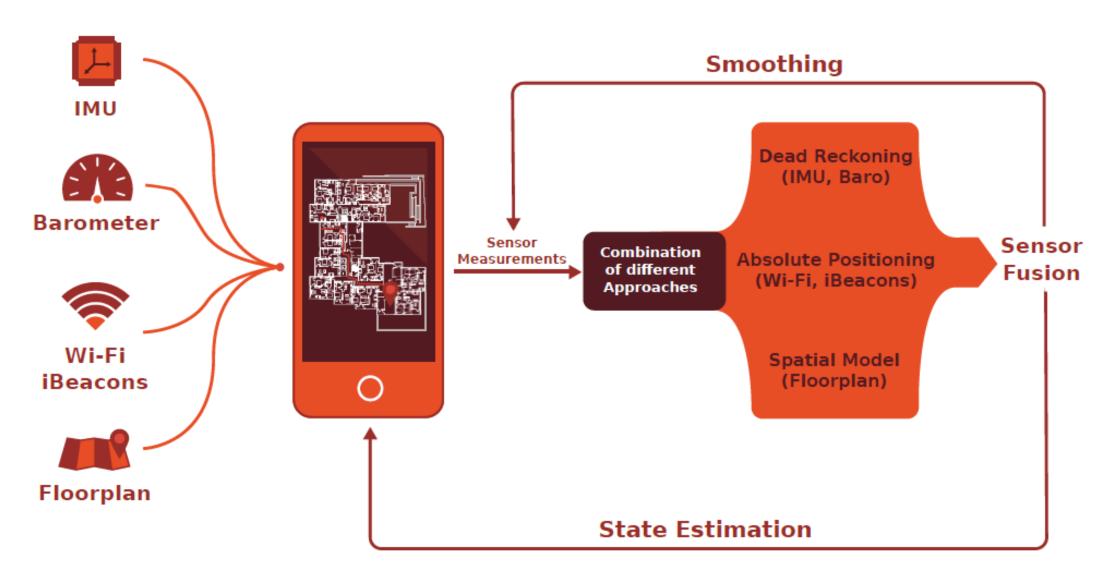
- Example particle filter: Catheter Tracking
 - ullet Transition $p(oldsymbol{q}_t | oldsymbol{q}_{t-1}, \langle oldsymbol{o}
 angle_{t-1})$
 - Along the way of the catheter probabilities shall be evenly distributed
 - Traditional transition without prior observations will avoid this
 - Integrating observations allows a formulation in the form

$$p(\boldsymbol{q}_t|\boldsymbol{q}_{t-1},\langle\boldsymbol{o}\rangle_{t-1}) \propto \frac{p(\boldsymbol{q}_t|\boldsymbol{q}_{t-1})}{\int\limits_{\boldsymbol{\epsilon}\in A}\int p(\boldsymbol{q}_t+\boldsymbol{\epsilon}|\boldsymbol{q}_{t-1})p(\boldsymbol{q}_{t-1}|\boldsymbol{o}_{t-1})d\boldsymbol{q}_{t-1}d\boldsymbol{\epsilon}}$$

- A describes a local area around q_t
- Denominator acts as smoothing of transition probability

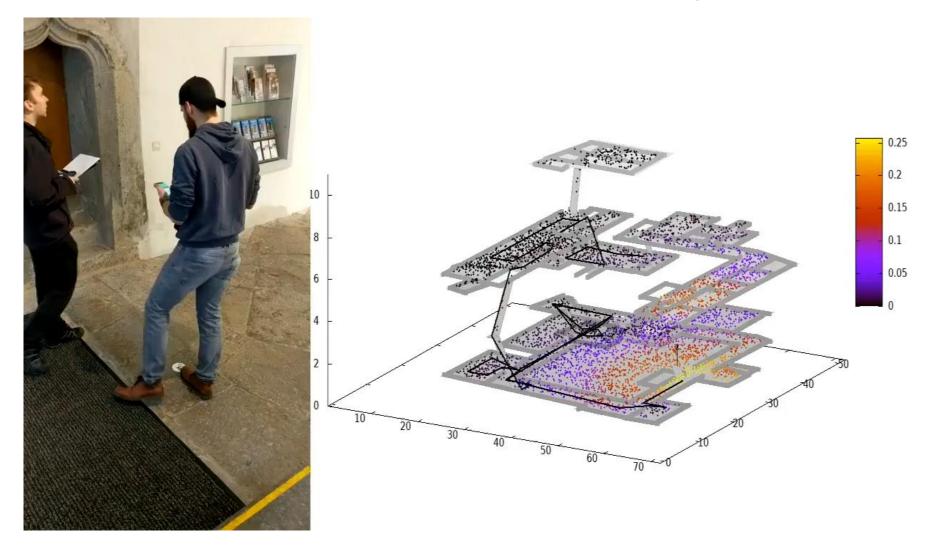


Particle Filter – Indoor Localization



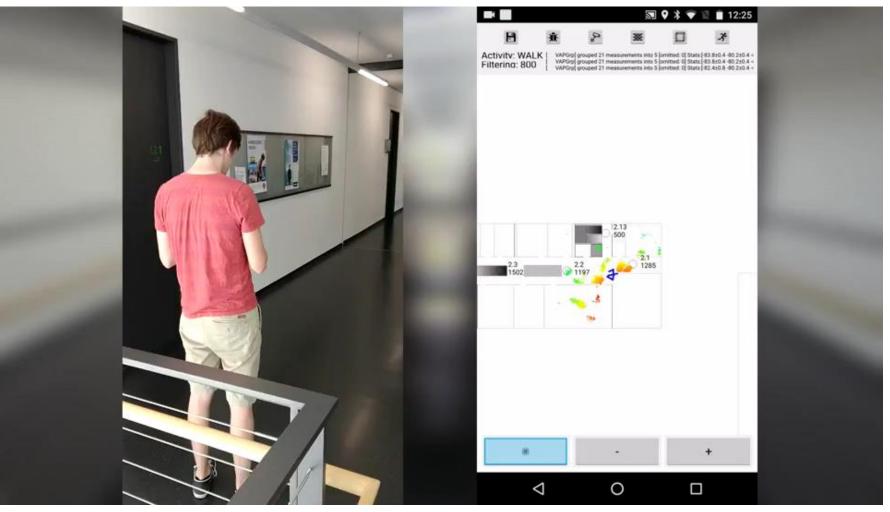
Particle Filter – Indoor Localization

► Example: Indoor Localization [Fet18], More Videos SimpleLoc YouTube



Particle Filter – Indoor Localization

► Example: Indoor Localization @ SHL, More Videos SimpleLoc YouTube

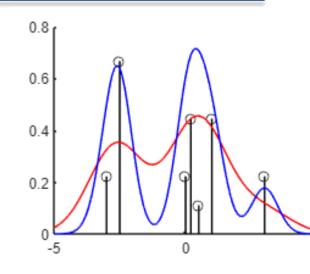


Particle Filtering Revisited

Approximate posterior by particle set (= set of weighted samples)

$$p(\langle \boldsymbol{q} \rangle_t \, | \, \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \, \langle \boldsymbol{o} \rangle_t) \approx \sum_i w_t^i \delta\left(\boldsymbol{s}_t^i - \boldsymbol{q}_t\right)$$

- Sample-based approximation has an obvious difficulty
 - Posterior is unknown
 - Hence sampling from posterior is impossible
- Solution: Importance Sampling
 - If we cannot sample from the desired posterior, then we sample from other density
 - We need to correct the error that we are making as a result of this
 - lacktriangle This other density is called *proposal density* or *importance density* and denoted by q



$$p(\langle \boldsymbol{q} \rangle_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t) = p(\boldsymbol{o}_t | \boldsymbol{q}_t) \int p(\boldsymbol{q}_t | \boldsymbol{q}_{t-1}) p(\boldsymbol{q}_{t-1} | \langle \boldsymbol{o} \rangle_{t-1}) d\boldsymbol{q}_{t-1}$$

- ullet When sampling from a proposal density q we need to compute the weights without knowing the posterior
 - Can be done this way [Thr05]

$$w_t^i \propto rac{p(\left\langle oldsymbol{s}^i
ight
angle_t | \left\langle oldsymbol{o}
ight
angle_t)}{q(\left\langle oldsymbol{s}^i
ight
angle_t | \left\langle oldsymbol{o}
ight
angle_t)}$$

- Luckily, we are not interested in the full sequence of states over time, but only in the current state
- Weighting can thus be reformulated similarly to the derivation of the recursive density estimate

$$w_t^i \propto w_{t-1}^i rac{p(\boldsymbol{o}_t|\boldsymbol{s}_t^i)p(\boldsymbol{s}_t^i|\boldsymbol{s}_{t-1}^i)}{q(\boldsymbol{s}_t^i|\boldsymbol{s}_{t-1}^i,\boldsymbol{o}_t)}$$

These are the weights for our sampled density

$$p(\boldsymbol{q}_t | \langle \boldsymbol{o} \rangle_t) \approx \sum_i w_t^i \delta\left(\boldsymbol{s}_t^i - \boldsymbol{q}_t\right)$$

■ This is the **Sequential Importance Sampling** (SIS) algorithm. These weights are optimal

At first glance, it looks as if we now have two unknown densities instead of one

$$w_t^i \propto w_{t-1}^i rac{p(\boldsymbol{o}_t|\boldsymbol{s}_t^i)p(\boldsymbol{s}_t^i|\boldsymbol{s}_{t-1}^i)}{q(\boldsymbol{s}_t^i|\boldsymbol{s}_{t-1}^i,\boldsymbol{o}_t)}$$

- It is up to you to use a suitable proposal density
 - Known from importance sampling: you can choose any proposal density you want (if they fulfill some basic properties)
 - Most common choice (Condensation algorithm does it that way) is

$$q(\boldsymbol{s}_t^i|\boldsymbol{s}_{t-1}^i,\boldsymbol{o}_t) = p(\boldsymbol{s}_t^i|\boldsymbol{s}_{t-1}^i)$$

 The transition model shall be the proposal density, because with that choice the weight update simplifies to

$$w_t^i \propto w_{t-1}^i p(\boldsymbol{o}_t | \boldsymbol{s}_t^i)$$

- Condensation algorithm
 - Sample from proposal density
 - Update weights

Condensation Algorithm

- 1. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from $S_{t ext{-}1}$ for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)

 The Condensation algorithm incorporates a resampling step to solve the degeneracy problem

- **Condensation Algorithm**
- 1. Create particle set S_0 for initial density $p(\mathbf{q}_0)$
- 2. Sample particles from S_{t-1} for new particle set S_t
- 3. Propagate each particle from S_t with state transition model
- 4. Evaluate each particle from S_t with observation model
- 5. Repeat from (2.)

- Performing without step (2.)
 - After few iterations weight of one particle will be close to 1.0
 - All other particle weights will be almost zero
 - This is the degeneracy problem
- Resampling of the particle set leads to new, equal weights
 - Degeneracy problem solved
- This is the **Sequential Importance Resampling** (SIR) filter
 - Transition prior used as proposal function
 - SIR filters are commonly known as Bootstrap Filter and Condensation Algorithm

- There are many more challenges in particle filtering
 - Sample Impoverishment: Transition step contains a deterministic and a stochastic part. Particles
 with same state will diversify during transition, but if noise variance is low, samples will not
 diversify enough. After some iterations particles will collapse into a single point in state space
 - Particle Filter Divergence: Poor tuning of filter, incorrect modeling assumptions, inconsistent measurements. Filter will diverge to "somewhere"
 - Proposal Density: Good proposal function is important. Transition as proposal is not always a good choice
 - Realtime Performance: There are many expensive components in a particle filter
 - Evaluation: Every particle must be evaluated
 - Transition: Every particle must be processed by transition model
 - Resampling: Particle set must be resampled. There are many different resampling strategies
 - Number of particles: Curse of dimensionality. Many parameters in state will lead to large particle sets
 - Good reading: [Elf21]

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Awards

- Winner of the "Indoor Localization Competition for Smartphone-Based Solutions" at the International Conference on Indoor Positioning and Indoor Navigation 2016 in Madrid, Spain
- Computer Science Award 2016 of the Fachbereichtag Informatik. Award for the master thesis Toni Fetzer "Smoothing and Prediction in Statistical Indoor Localization"
- Best-Paper Award of the International Conference on Indoor Positioning and Indoor Navigation 2015 (Frank Ebner, Toni Fetzer, Lukas Koeping, Marcin Grzegorzek, Frank Deinzer): Award for the paper "Multi Sensor 3D Indoor Localization".
- Innovationspreis 2008 der Gesellschaft für Informatik e.V. (Frank Deinzer, Esther Platzer): Auszeichnung für die Arbeit "Erstellung von 4-D-Angiogrammen in der interventionellen Radiologie"
- Scientific Award 2007 der BMW Group (1. Platz): Auszeichnung der Diplomarbeit Esther Platzer: "Visualierung von Blutfluss in 3-D-Datensätzen aus 2-D-Angiogrammen"
- DAGM-Preis 2001 (Frank Deinzer, Joachim Denzler, Heinrich Niemann): Auszeichnung für die Arbeit "On Fusion of Multiple Views for Active Object Recognition"

