

With the prior information regarding that the target has been detected affect the state distribution?

$$\begin{aligned} p(x|\delta) &= \frac{p(\delta|x)p(x)}{p(\delta)} = \frac{p(\delta|x)p(x)}{\int p_{\delta}(x)p(x)dx} \\ &= \frac{p_{\delta}(x)p(x)}{\int p_{\delta}(x)p(x)dx} \end{aligned}$$

The question is whether this changes the distribution.

Assuming $p_{\delta}(x)$ constant, we get that

$$p(x|\delta) = \frac{p_{\delta} p(x)}{p_{\delta} \int p(x)dx} = p(x)$$

however, by assuming that $p_{\delta}(x)$ varying, we get that

$$p(x|\delta) = \frac{p_{\delta}(x)p(x)}{\int p_{\delta}(x)p(x)dx} \neq p(x)$$

assuming that $\int p_{\delta}(x) \neq p_{\delta}(x)p(x)$

in other words, prior information of existence changes the distribution

a) Sensor with N cells and a likelihood of false alarm given as P_{FA}

Assuming that there are M_N sensor cells being detected used.

number of

The false alarms is given us a binomial distribution:

$$P(\varphi_n = \varphi) = \binom{M_N}{\varphi} P_{FA}^{\varphi} (1-P_{FA})^{M_N-\varphi}$$

where φ with $0 \leq \varphi \leq M_N \leq N$

b) Assuming a Poisson-distribution with $\lambda = \frac{1}{V_n} M_N P_{FA}$

Supposed to use that

$$P_{\delta}\{a_n | \gamma_{1:n}\} \propto \begin{cases} (1-p_{\delta})^{M_N} \frac{\lambda^{M_N}}{M_N!}, & a_n = 0 \\ \frac{p_{\delta}^n \lambda^{a_n}}{C(a_n)} & , a_n > 0 \end{cases}$$

I will be relying a bit on the lectures, as I would otherwise have been quite high to come up with this method.
we know that

$$z_u^{au} = \int dz (z_u^{au} | x_u) p_{\text{prior}}(x_u) dz$$

Under gaussian-linear assumptions, we have that

$$z_u^{au} \sim N(z_u^{au}; z_{u|u-1}, s_u)$$

This gives us that (by also abusing the proportionality)

$$\Pr\{a_u | z_{1:n}\} \propto \begin{cases} (1-p_D) \mu(m_u) & , a_u = 0 \\ \frac{p_D \mu(m_{u-1})}{m_u(z_u)} N(z_u^{au}; z_{u|u-1}, s_u) & , a_u > 0 \end{cases}$$

By using that this is a zero poisson-distribution, we have that

$$\mu(\lambda) = \frac{\lambda^u}{u!} \exp(-\lambda)$$

which gives that

$$\Pr\{a_u | z_{1:n}\} \propto \begin{cases} (1-p_D) \frac{1^{m_u} \exp(-\lambda)}{m_u!} & , a_u = 0 \\ \frac{p_D \lambda^{m_{u-1}} \exp(-\lambda)}{(m_{u-1})! m_u} N(z_u^{au}; z_{u|u-1}, s_u) & , a_u > 0 \end{cases}$$

$$\propto \begin{cases} (1-p_D) \lambda^{m_u} & , a_u = 0 \\ p_D N(z_u^{au}; z_{u|u-1}, s_u) & , a_u > 0 \end{cases}$$

c) since we are to use real distribution from a), I assume that we are to use the binomial-distribution

This gives that

$$\mu(m_u) = \binom{m_u}{m_u} p_{\text{TA}}^{m_u} (1-p_{\text{TA}})^{m_u-m_u}$$

$$\Pr\{a_u | z_{1:n}\} \propto \begin{cases} (1-p_D) \binom{m_u}{m_u} p_{\text{TA}}^{m_u} (1-p_{\text{TA}})^{m_u-m_u} & , a_u = 0 \\ \frac{p_D \binom{m_u}{m_{u-1}} p_{\text{TA}}^{m_{u-1}} (1-p_{\text{TA}})^{m_u-m_u+1}}{m_u (z_u)} N(z_u^{au}; z_{u|u-1}, s_u) & , a_u > 0 \end{cases}$$

By abusing the proportionality sign, this gives us that

$$\therefore \text{cc} \left\{ (1-P_D) \frac{m_k P_{FA}}{1-P_{FA}} \binom{m_k}{m_k} \cdot \frac{1}{\binom{m_k}{m_k-1}}, m_k \geq 0 \right\}$$

$$P_D N(z_k^{m_k}; z_{k|m_k-1}^*, s_k) , m_k \geq 0$$

$$\binom{m_k}{m_k} \cdot \frac{1}{\binom{m_k}{m_k-1}} = \frac{m_k!}{m_k!(m_k-m_k)!} \cdot \frac{(m_k-1)! (m_k-m_k+1)!}{m_k!}$$

$$= \frac{(m_k-m_k+1)!}{m_k}$$

$$\Rightarrow \text{cc} \left\{ (1-P_D) P_{FA} m_k \frac{(m_k-m_k+1)!}{m_k (1-P_{FA})}, m_k \geq 0 \right\}$$

$$P_D N(z_k^{m_k}; z_{k|m_k-1}^*, s_k) , m_k \geq 0$$

I cannot see how one should simplify the top expression further.

ii) assuming that m_k is large and P_{FA} small. This gives that

$$\text{cc} \left\{ \Pr \{ \text{at least } k \text{ hits} \} \right\} \underset{\substack{P_{FA} \rightarrow 0 \\ m_k \rightarrow \infty}}{\approx} \frac{(1-P_D)^k}{\underbrace{\frac{m_k-1}{1-P_{FA}}}_k} = (1-P_D)^k$$

$$\rightarrow 1$$

As expected due to the properties of the poisson-distribution and the binomial-distribution, the approximation seems good.

3a) The linear-gaussian assumption assumes that all of the distributions are gaussian, such that the ~~posterior~~ posterior is ~~a~~ gaussian.

The PDAT tries to combine all of the measurement-models into a single distribution by mixing them. This is done to save computational power by only keeping track of one mixture, but also allows a weighting of the measurement models. Under the simple-target tracking, it is impossible to know exactly which measurement has hit upon the target, due to cluttering and/or mis detections.

By weighting the measurements, one could quantify how likely it is that ~~one~~ a measurement originates from the target.

- b) ~~is a single target tracking algorithm that~~
JPDA ~~tries to track a target using when there is no information whether the target exists or not. In other words, the JPDA ~~is~~ is PDA (7) with the added estimation if the target is present or not.~~
Since PDA ~~&~~ JPDA, there are no problems that PDA solves that JPDA cannot.

- c) The JPDA will have a larger state-estimation compared to the PDA. It cannot just implicitly assume that a target exists. It has to find and maintain such a likelihood as well.

```
def get_mean(self) -> ndarray:  
    """  
    Return the mean of the gaussian mixture  
  
    Hint: use what you did in mixturereductin.py assignment 4  
  
    Returns:  
        mean (ndarray): the mean  
    """  
  
    # I assume that the system should mixture self.gaussians  
    means = np.zeros_like(self.weights)  
  
    # Extract all of the data from the gaussians  
    i = 0  
    for gaussian in self.gaussians:  
        means[i] = gaussian.mean  
        i += 1  
  
    # Multiply by the weights of the corresponding gaussians  
    mean = self.weights.T @ means  
  
    #mean = solution.gaussmix.GaussianMuxture.get_mean(self)  
    return mean
```

```
def get_cov(self) -> ndarray:  
    """  
    Return the covariance of the gaussian mixture  
  
    Hint: use what you did in mixturereductin.py assignment 4  
  
    Returns:  
        cov (ndarray): the covariance  
    """  
  
    mean = self.get_mean()  
    #diffs = means - mean[np.newaxis]  
  
    cov_shape = np.shape(self.gaussians[0].cov)  
    covs = np.zeros_like(cov_shape)  
    #for i in range(len(self.gaussians)):  
    #    covs[i*cov_shape[0]:i]  
  
    #cov_internal = np.average(covs, axis=0, weights=self.weights)  
    #cov = 0  
  
    # I was unable to actually solve this in task 4, and still haven't  
    # understood how one should solve this  
  
    cov = solution.gaussmix.GaussianMuxture.get_cov(self)  
    return cov
```

```
def reduce(self) -> MultiVarGaussian:  
    """  
    Reduce the gaussian mixture to a multivariate gaussian  
    Hint: you can use self.get_mean and self.get_cov  
  
    Returns:  
        reduction (MultiVarGaussian): the reduction  
    """  
  
    mean = self.get_mean()  
    cov = self.get_cov()  
  
    reduction = np.random.multivariate_normal(mean, cov)  
  
    #reduction = solution.gaussmix.GaussianMuxture.reduce(self)  
    return reduction
```

```
def predict_state(  
    self,  
    state_upd_prev_gauss: MultiVarGaussian,  
    Ts: float  
) -> MultiVarGaussian:  
    """  
    Prediction step  
    Hint: use self.ekf  
  
    Args:  
        state_upd_prev_gauss (MultiVarGaussian): previous update gaussian  
        Ts (float): timestep  
  
    Returns:  
        state_pred_gauss (MultiVarGaussian): predicted state gaussian  
    """  
  
    state_pred_gauss = self.ekf.predict_state(  
        state_upd_prev_gauss=state_upd_prev_gauss,  
        Ts=Ts)  
  
    # state_pred_gauss = solution.pdaf.PDAF.predict_state(  
    #     self, state_upd_prev_gauss, Ts)  
    return state_pred_gauss
```

```
def predict_measurement(
    self,
    state_pred_gauss: MultiVarGaussian
) -> MultiVarGaussian:
"""

Measurement prediction step
Hint: use self.ekf

Args:
    state_pred_gauss (MultiVarGaussian): predicted state gaussian

Returns:
    z_pred_gauss (MultiVarGaussian): predicted measurement gaussian
"""

z_pred_gauss = self.ekf.predict_measurement(state_pred_gauss=state_pred_gauss)

# z_pred_gauss = solution.pdaf.PDAF.predict_measurement(
#     self, state_pred_gauss)
return z_pred_gauss
```

```
def gate(
    self,
    z_pred_gauss: MultiVarGaussian,
    measurements: Sequence[ndarray]
) -> ndarray:
"""

Gate the incoming measurements. That is remove the measurements
that have a mahalanobis distance higher than a certain threshold.

Hint: use z_pred_gauss.mahalanobis_distance_sq and self.gate_size_sq

Args:
    z_pred_gauss (MultiVarGaussian): predicted measurement gaussian
    measurements (Sequence[ndarray]): sequence of measurements

Returns:
    gated_measurements (ndarray[:,2]): array of accepted measurements
"""

gated_measurements = np.empty_like([measurements[0]])
is_initialized = False

for m in measurements:
    if z_pred_gauss.mahalanobis_distance_sq(m) > self.gate_size_sq:
        continue
    if is_initialized:
        gated_measurements = np.append(gated_measurements, np.array(m))
        continue
    gated_measurements[:] = np.array(m)
    is_initialized = True

# gated_measurements = solution.pdaf.PDAF.gate(
#     self, z_pred_gauss, measurements)
return gated_measurements
```

```
def get_association_prob(  
    self,  
    z_pred_gauss: MultiVarGaussian,  
    gated_measurements: ndarray  
) -> ndarray:  
    """  
    Finds the association probabilities.  
  
    associations_probs[0]: prob that no association is correct  
    associations_probs[1]: prob that gated_measurements[0] is correct  
    associations_probs[2]: prob that gated_measurements[1] is correct  
    ...  
  
    the sum of associations_probs should be 1  
  
    Args:  
        z_pred_gauss (MultiVarGaussian): predicted measurement gaussian  
        gated_measurements (ndarray[:,2]): array of accepted measurements  
  
    Returns:  
        associations_probs (ndarray[:]): the association probabilities  
    """  
    P_D = self.detection_prob  
    V = 1 / self.clutter_density  
  
    m = len(gated_measurements)  
    associations_probs = np.zeros((m,1))  
    associations_probs = np.append(associations_probs, np.zeros((1,1)))  
  
    # Calculating for i == 0  
    associations_probs[0] = m/V * (1 - P_D)  
  
    # Calculating for i > 0  
    for i in range(1, m):  
        I_i = z_pred_gauss.pdf(gated_measurements[i])  
        associations_probs[i-1] = P_D*I_i  
  
    associations_probs /= associations_probs.sum()  
  
    # associations_probs = solution.pdaf.PDAF.get_association_prob(  
    #     self, z_pred_gauss, gated_measurements)  
    return associations_probs
```

```
def get_cond_update_gaussians(
    self,
    state_pred_gauss: MultiVarGaussian,
    z_pred_gauss: MultiVarGaussian,
    gated_measurements: ndarray
) -> Sequence[MultiVarGaussian]:
```

```
# Allocating memory
n = len(gated_measurements)
update_gaussians = np.arange(n+1, dtype=MultiVarGaussian)

# Get the associated probabilities
associated_probs = self.get_association_prob(z_pred_gauss, gated_measurements)

# Iterate over the associated probabilities and calculate the posteriori state given
# that the associated measurement is correct
for i in range(len(associated_probs)):
    gaussian_mixture_class = GaussianMixture(
        weights=np.array([associated_probs[i]]),
        gaussians=np.array([state_pred_gauss]))
    update_gaussians[i] = gaussian_mixture_class.reduce()

# Something is incorrect with this code, since it will not reduce properly
# for all

update_gaussians = solution.pdaf.PDAF.get_cond_update_gaussians(
    self, state_pred_gauss, z_pred_gauss, gated_measurements)
return update_gaussians
```

```
def update(
    self,
    state_pred_gauss: MultiVarGaussian,
    z_pred_gauss: MultiVarGaussian,
    measurements: Sequence[ndarray]
) ->MultiVarGaussian:
"""
Perform the update step of the PDA filter

Args:
    state_pred_gauss (MultiVarGaussian): predicted state gaussian
    z_pred_gauss (MultiVarGaussian): predicted measurement gaussian
    measurements (Sequence[ndarray]): sequence of measurements

Returns:
    state_upd_gauss (MultiVarGaussian): updated state gaussian
"""

# Gate the measurements
gated_measurements = self.gate(
    z_pred_gauss=z_pred_gauss,
    measurements=measurements)

# Get the associated probabilities
associated_probs = self.get_association_prob(
    z_pred_gauss=z_pred_gauss,
    gated_measurements=gated_measurements)

# Get the conditional associated states
cond_states = self.get_cond_update_gaussians(
    state_pred_gauss=state_pred_gauss,
    z_pred_gauss=z_pred_gauss,
    gated_measurements=gated_measurements)

gaussian_mixture_class = GaussianMixture(
    weights=associated_probs,
    gaussians=cond_states)
state_upd_gauss = gaussian_mixture_class.reduce()

# state_upd_gauss is here just an array of float64
# and not a gaussian... Just fuck python! C++ all the way (except for plotting)

state_upd_gauss = solution.pdaf.PDAF.update(
    self, state_pred_gauss, z_pred_gauss, measurements)
return state_upd_gauss
```

```
def step_with_info(
    self,
    state_upd_prev_gauss: MultiVarGaussian,
    measurements: Sequence[ndarray],
    Ts: float
) -> Tuple[MultiVarGaussian,
            MultiVarGaussian,
            MultiVarGaussian]:
"""

Perform a full step and return usefull info

Hint: you should not need to write any new code here,
just use the methods you have implemented

Args:
    state_upd_prev_gauss (MultiVarGaussian): previous updated gaussian
    measurements (Sequence[ndarray]): sequence of measurements
    Ts (float): timestep

Returns:
    state_pred_gauss (MultiVarGaussian): predicted state gaussian
    z_pred_gauss (MultiVarGaussian): predicted measurement gaussian
    state_upd_gauss (MultiVarGaussian): updated state gaussian
"""

state_pred_gauss = self.predict_state(state_upd_prev_gauss=state_upd_prev_gauss, Ts=Ts)
z_pred_gauss = self.predict_measurement(state_pred_gauss=state_pred_gauss)
state_upd_gauss = self.update(state_pred_gauss=state_pred_gauss, z_pred_gauss=z_pred_gauss, measurements=measurements)

# state_pred_gauss, z_pred_gauss, state_upd_gauss = solution.pdaf.PDAF.step_with_info(
#     self, state_upd_prev_gauss, measurements, Ts)
return state_pred_gauss, z_pred_gauss, state_upd_gauss
```

....

Been raging a bit too hard on python and numpy, so my bloodpressure is a bit too high right now...

Also depressed, such that I have no motivation on tuning the filter well...

Perfect time to tune in other words!

I will try to rather get a feeling about the response, instead of having a filter that is "perfect"....

Once my blood starts cooling down, perhaps I try to get a more optimized filter, but cannot guarantee that!

`sigma_a` and `sigma_z` is related to the EKF

`sigma_z = R` and is the noise belonging to the measurements. By increasing this, we expect that the measurements will have a larger noise, and thus trust the model more. For the PDA, this means that values that are further away from the measurement is more related to noise in measurements instead of the actual process. Increasing this should in fact reduce number of samples that will actively be considered

`sigma_z = Q` is the noise related to the model. By increasing this, we say that the process can change rapidly. This allows us to increase the likelihood for measurements further away from the priori estimate. Thus, considering more estimates

Clutter density is number of measurements per m^2 . Increasing this allows the system to generate more clutter. I cannot identify if this exact value is reasonable or not, however the clutter density will vary depending on the accuracy of the system as well as of the desired scan-velocity.

For example would a radar have lower clutter intensity with a lower rotation speed, with the downside that it updates relatively infrequently. Increasing the rotation speed allows it to scan the area quicker, however will allow less time for the sensor cells to acquire and identify a possible signal.

Detection probability should be in the range [0.5, 1) Lower bounded by 0.5 since that value means that we are just guessing yes/no if a detection has been made.

Note written couple of days later:

I should have tuned this a lot better. The difficult thing is to get the system to follow the correct state without diverging too rapidly. My experimentation showed that the system would either accept too many measurements and therefore diverge, or that it would not value the real system's measurements and thus diverge. I did experience that the algortihm was able to follow the track somewhat, however the filter became overconfident and started diverging at iteration 136. However, when running the same value on my desktop (right when I am writing this shit), the system diverged at iteration number 7 to 8. Idk why my laptop and desktop gave different results....

I honestly started by trying to find the values that made the system stay within a reasonable NEES, however that didn't quite work out.

After that, I tried to just tune after analysing the track... It went as well as one could expect.

In other words, fuck my life!

.....

