# Sensorfusion assignment 4

#### Bjørn Kåre Sæbø

#### Task 1

## A)

```
Implemented based on 6.19 - 6.22 in book
"""Calculate the first two moments of a Gaussian mixture"""

# mean
mean_bar = np.average(mean, weights=w, axis=0)

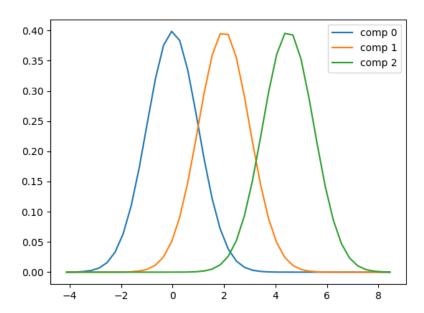
# covariance
# # internal covariance
cov_int = np.average(cov, weights=w, axis=0)

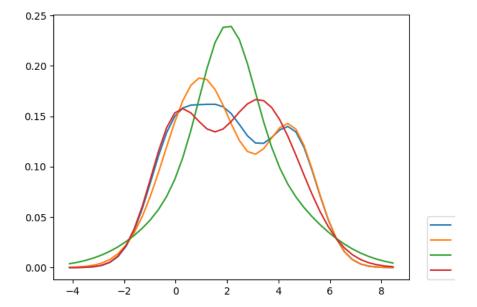
# * spread of means
# Optional calc: mean_diff =
mean_diff = np.diagonal((mean - mean_bar)@(mean - mean_bar).T)
cov_ext = np.average(mean_diff, weights=w, axis=0)

# # total covariance
cov_bar = cov_int + cov_ext
return mean_bar, cov_bar
```

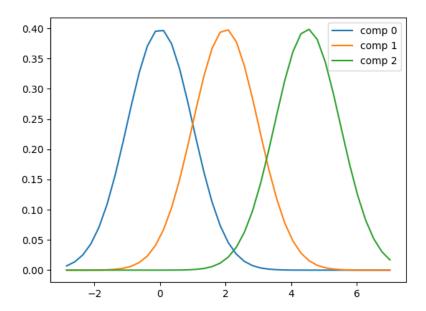
B)

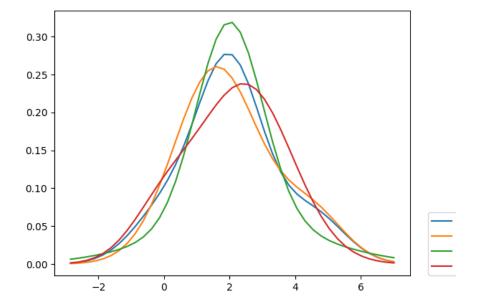
I) Would merge the outer distributions to get a wider distribution with mean in the middle of the three



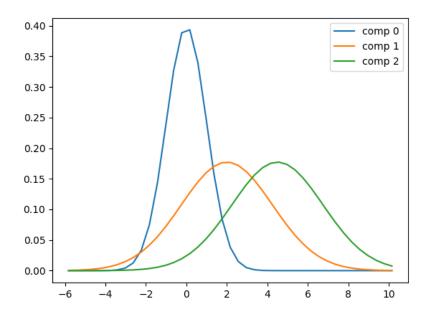


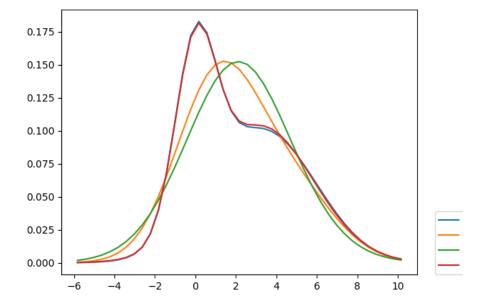
II) Doesn't make too much difference, as the weighting ensures that the middle distribution dominates the result anyway



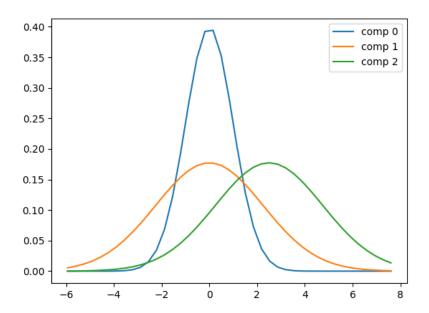


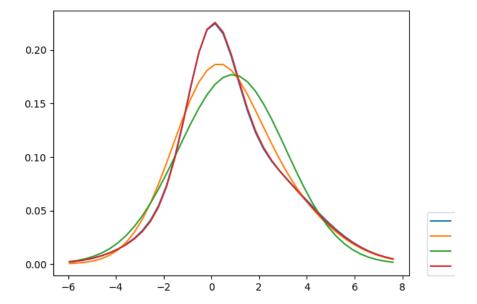
III) Would merge the two wide rightmost distributions, as they contribute much of the same info



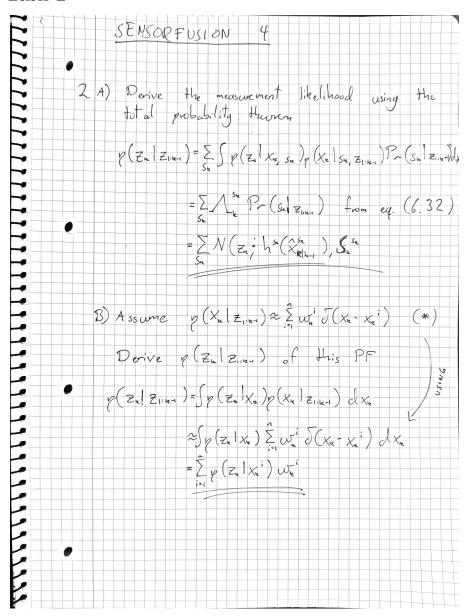


IV) Would merge the distributions with mean 0, and get a distribution with mean 0 and variance somewhere between the two originals





### Task 2



### Task 3 & 4

Spent a lot of time trying to get this to work, eventually gave up. Not available on friday, so can't bother the student assistant any more than I already have. Including code from IMM.py for reference

```
class IMM(Generic[MT]):
    # The M filters the IMM relies on
   filters: List[StateEstimator[MT]]
    # the transition matrix. PI[i, j] = probability of going from model i to j: shape (M, M.
   PI: np.ndarray
    # init mode probabilities if none is given
    initial_mode_probabilities: Optional[np.ndarray] = None
    def __post_init__(self):
        # This have to be satisfied!
        if not np.allclose(self.PI.sum(axis=1), 1):
            raise ValueError("The rows of the transition matrix PI must sum to 1.")
        # Nice to have a reasonable initial mode probability
        if self.initial_mode_probabilities is None:
            eigvals, eigvecs = linalg.eig(self.PI)
            self.initial_mode_probabilities = eigvecs[:, eigvals.argmax()]
            self.initial_mode_probabilities = (
                self.initial_mode_probabilities / self.initial_mode_probabilities.sum()
    def mix_probabilities(
        self,
        immstate: MixtureParameters[MT],
        # sampling time
        Ts: float,
    ) -> Tuple[
        np.ndarray, np.ndarray
   ]: # predicted_mode_probabilities, mix_probabilities: shapes = ((M, (M ,M))).
        # mix_probabilities[s] is the mixture weights for mode s
        """Calculate the predicted mode probability and the mixing probabilities."""
        predicted mode probabilities, mix probabilities = (
            discretebayes.discrete_bayes(immstate.weights,self.PI)
        # Optional assertions for debugging
        assert np.all(np.isfinite(predicted_mode_probabilities))
        assert np.all(np.isfinite(mix_probabilities))
        assert np.allclose(mix_probabilities.sum(axis=1), 1)
        return predicted_mode_probabilities, mix_probabilities
    def mix_states(
        self.
        immstate: MixtureParameters[MT],
```

```
# the mixing probabilities: shape=(M, M)
   mix_probabilities: np.ndarray,
) -> List[MT]:
   mixed_states = [f1.reduce_mixture(MixtureParameters(mp, immstate.components))
                    for fl, mp in zip(self.filters, mix_probabilities)]
    return mixed_states
def mode_matched_prediction(
    self,
   mode_states: List[MT],
    # The sampling time
   Ts: float,
) -> List[MT]:
   pred = []
   for m in mode states:
        for fl in self.filters:
            pred.append(fl.predict(m, Ts))
   return np.array(pred)
def predict(
    self,
    immstate: MixtureParameters[MT],
    # sampling time
   Ts: float,
) -> MixtureParameters[MT]:
    Predict the immstate Ts time units ahead approximating the mixture step.
    Ie. Predict mode probabilities, condition states on predicted mode,
    appoximate resulting state distribution as Gaussian for each mode, then predict each
    HHHH
    predicted_mode_probability, mixing_probability = self.mix_probabilities(immstate, Table 1)
   mixed_mode_states: List[MT] = self.mix_states(immstate,mixing_probability)
    predicted_mode_states = self.mode_matched_prediction(mixed_mode_states, Ts)
   predicted_immstate = MixtureParameters(
        predicted_mode_probability, predicted_mode_states
   return predicted_immstate
def mode_matched_update(
```

```
self,
    z: np.ndarray,
    immstate: MixtureParameters[MT],
    sensor_state: Optional[Dict[str, Any]] = None,
) -> List[MT]:
    """Update each mode in immstate with z in sensor_state."""
    updated_state = np.array([fl.update(z, m, sensor_state) for fl, m in zip(self.filter
    return updated_state
def update_mode_probabilities(
    self,
    z: np.ndarray,
    immstate: MixtureParameters[MT],
    sensor_state: Dict[str, Any] = None,
) -> np.ndarray:
    """Calculate the mode probabilities in immstate updated with z in sensor_state"""
    mode_loglikelihood = np.array(
        [filt.loglikelihood(z, comp, sensor_state) \
         for filt, comp in zip(self.filters, immstate.components)])
    # potential intermediate step logjoint =
   predicted_mode_probabilities = immstate.weights # shape (M,1)
   normalization = \
        np.sum(np.exp(mode_loglikelihood) * predicted_mode_probabilities) # scalar floa
    log_pred_mode_probs = np.log(predicted_mode_probabilities) # shape (M,1)
    log_norm = np.log(normalization) # scalar float
    # (6.33)
    # mode_loglikelihood * predicted_mode_probabilities / normalization
    log updated mode probs = \
        mode_loglikelihood + log_pred_mode_probs - log_norm
    updated_mode_probabilities = np.exp(log_updated_mode_probs)
    # Optional debuging
    assert np.all(np.isfinite(updated_mode_probabilities))
    assert np.allclose(np.sum(updated_mode_probabilities), 1)
    return updated_mode_probabilities
def update(
    self,
    z: np.ndarray,
```

```
immstate: MixtureParameters[MT],
    sensor_state: Dict[str, Any] = None,
) -> MixtureParameters[MT]:
    """Update the immstate with z in sensor_state."""
    updated_weights = self.update_mode_probabilities(z, immstate)
    updated_states = self.mode_matched_update(z, immstate)
    updated_immstate = MixtureParameters(updated_weights, updated_states)
    return updated_immstate
def step(
   self,
   z,
    immstate: MixtureParameters[MT],
   Ts: float,
    sensor_state: Dict[str, Any] = None,
) -> MixtureParameters[MT]:
    """Predict immstate with Ts time units followed by updating it with z in sensor_sta
    predicted_immstate = self.predict(immstate, Ts)
    updated_immstate = self.update(z, immstate, Ts)
   return updated_immstate
def loglikelihood( # Postponed until required
    self,
    z: np.ndarray,
    immstate: MixtureParameters,
    sensor_state: Dict[str, Any] = None,
) -> float:
    # THIS IS ONLY NEEDED FOR IMM-PDA. You can therefore wait if you prefer.
   mode_conditioned_ll = None # TODO in for IMM-PDA
    11 = None # TODO
    return 11
def reduce_mixture(
    self, immstate_mixture: MixtureParameters[MixtureParameters[MT]]
) -> MixtureParameters[MT]:
    """Approximate a mixture of immstates as a single immstate"""
```

```
# extract probabilities as array
    weights = immstate_mixture.weights
    component_conditioned_mode_prob = np.array(
        [c.weights.ravel() for c in immstate_mixture.components]
    # flip conditioning order with Bayes
    mode_prob, mode_conditioned_component_prob = discretebayes.discrete_bayes(weights, o
    # Hint list_a of lists_b to list_b of lists_a: zip(*immstate_mixture.components)
   mode_states = None # TODO:
    immstate_reduced = MixtureParameters(mode_prob, mode_states)
   return immstate_reduced
def estimate(self, immstate: MixtureParameters[MT]) -> GaussParams:
    """Calculate a state estimate with its covariance from immstate"""
    # ! You can assume all the modes have the same reduce and estimate function
    # ! and use eg. self.filters[0] functionality
    data_reduced = [gaussian_mixture_moments(immstate.weights, comp.mean, comp.cov) for
    estimate = self.filters[0].estimate(data_reduced)
   return estimate
```

#### Task 5

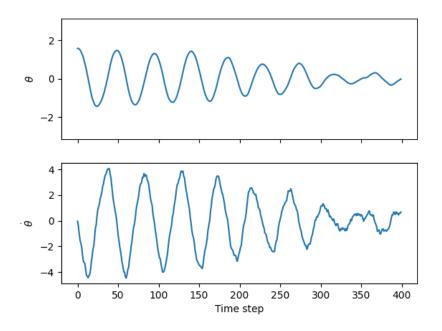
#### A) Finished code

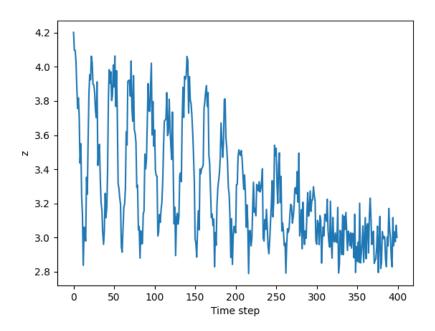
```
for k in range(K):
    print(f"k = {k}")
    # weight update
    for n in range(N):
        dz = Z[k] - h(px[n], Ld, 1, L1)
        w[n] = PF_measurement_distribution.pdf(dz)
    w += eps
    w = w / np.sum(w)

# resample
# TODO: some pre calculations?
# N_eff = 1 / np.sum(w**2)
# if N_eff <= N/2:
# resample using algorithm 3 p. 90
cumweights = np.cumsum(w)</pre>
```

```
indicesout = np.zeros((N, 1))
noise = rng.random((1,1)) / N
i = 0
for n in range(N):
    # find a particle 'i' to pick
    # algorithm in the book, but there are other options as well
    u_n = n/N + noise
    while u_n > cumweights[i]:
       i += 1
    pxn[n] = px[i]
# indicesout = indicesout[np.randperm(np.size(indicesout))]
rng.shuffle(pxn,axis=0)
w.fill(1/N)
# else:
# pxn = px[:]
N_{eff} = 1 / np.sum(w**2)
# trajecory sample prediction
for n in range(n):
    vkn = PF_dynamic_distribution.rvs()
    px[n] = pendulum_dynamics_discrete(pxn[n], vkn, Ts, a)
```

This worked decently enough with 100 filter, but became very noisy towards the end, when the pendulum slowed





### B)

Larger  $L_l$  seems to give better results, which makes some sense, based on intuition from the bearings-only example in the lecures.

# C)

PF seems like a good alternative to EKFs when there is a lot of nonlinearity in process models and measurement models