

## 02471 Machine Learning for Signal Processing

# Assignment 3

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## 1 Problem 3.1 Sparse signal representations

#### 1.1 Problem 3.1.1

A multitone signal is sparse in the DCT domain (we can see that by looking at the DCT formula (week 7 slide 22). Then, we can estimate the signal from a few samples by taking this sparsity into account. The goal is to find an acceptable estimation of the original signal in the sparse domain that has as few non-zero components as possible (the smallest components must be put down to zero). This is possible by minimizing the least square error under some constraints which is equivalent to adding a regularization term on the least square error cost function. The LASSO cost function (equation (9.6) in ML) uses the L1 norm in the regularization term. By minimizing this cost function, we can derive equation (9.13) in ML which shows that if an estimated value of the signal using LS estimate is too small compared to lambda it will be put down to zero in the LASSO estimation. Lambda controls the amount of non zero-elements. We can use other norms such as the L0 norm which is the sparsest norm but the L1 norm is the computationally most efficient norm as it is the most sparse true norm, and is convex. In practice, the cost function could be minimized using the gradient descent as we have seen in the IST algorithm.

#### 1.2 Problem 3.1.2

I decided to use the LASSO cost function and IST. Both to solve this problem as compare to each other. I used the build-in function lasso in Python and IST algorithm. I choose  $\lambda = 0.01$  as it gives an acceptable number of non-zero components. The estimated signal is displayed figure 1 both in time and DCT domain. The estimated parameters of the multitone signal are given below: (As too much values I only print a and m with first 4 respective)

- K status: 36
- $a_i$  values:  $\begin{bmatrix} 2.84376962e-01 & 7.38536161e-01 & 1.36582746e-02 & 8.04106128e-01 \end{bmatrix}$
- $m_j$  values: [ 4 9 22 24]

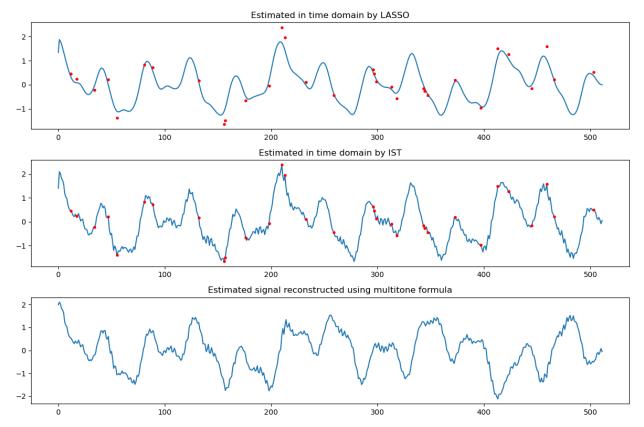


Figure 1: Sampled signal.

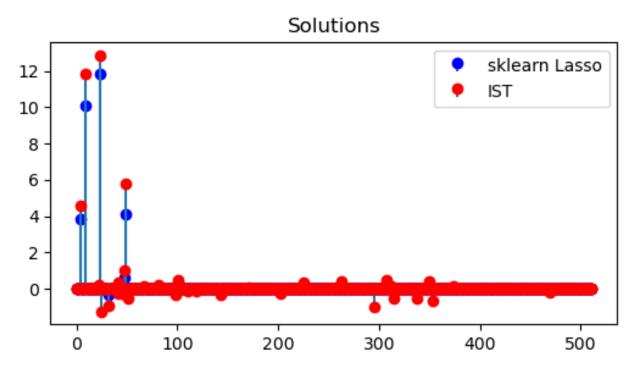


Figure 3: DCT domain.

### 2 Problem 3.2 Bayesian inference and the EM algorithm

#### 2.1 Problem 3.2.1

In this optimization problem, the two terms can be interpreted as the likelihood, prior, mean square error, and regularization.

#### Roles of the Objective Terms:

- $\frac{1}{2\sigma_n^2}||y X\theta||_2^2$ :
  - The perspectives of likelihood and mean squared error:
    - \* Quantifies the error between the model's predictions  $X\theta$  and the observed data y.
    - \* Appears as the likelihood, minimizing this error encourages in order to maximize the likelihood of observed data y given the model.
- $\bullet \ \frac{1}{2\sigma_{\theta}^2}||\theta||_2^2$ :
  - Prior Knowledge and Regularization Perspective:
    - \* Represents the negative log of the prior probability distribution for model parameters  $\theta$ .
    - \* Acts as a regularization term, controlling the size of parameter  $\theta$ , preventing over-reliance on parameters, avoiding overfitting, and enhancing model generalization.

#### Influence of Parameters $\sigma_{\eta}$ and $\sigma_{\theta}$ :

- $\sigma_{\eta}$ (Standard deviation of observation errors):
  - A larger  $\sigma_{\eta}$  indicates higher noise levels in observed data y, leading to a greater reliance on the regularization term, reducing overfitting.
  - A smaller  $\sigma_{\eta}$  indicates lower noise levels in observed data y, resulting in a greater reliance on the MSE term, increasing the fit to observed data.
- $\sigma_{\theta}$  (Standard deviation of parameter  $\theta$ ):
  - A larger  $\sigma_{\theta}$  implies less restriction on  $\theta$ , allowing wider ranges for parameter values.  $\theta$ .
  - A smaller  $\sigma_{\theta}$  implies stricter constraints on  $\theta$ , leading to more restricted parameter values.

These parameter choices balance the model's fit to observed data and control over parameter sizes. Adjusting  $\sigma_{\eta}$  and  $\sigma_{\theta}$  values can alter the model's performance to adapt to varying noise levels and complexities.



#### 2.2 Problem 3.2.2

As we know the density function,

$$p(z|a,b) = \frac{b^a}{\Gamma(a)} z^{-a-1} exp(-\frac{b}{z})$$

we take log of the density function  $\ln p(z|a,b)$ ,

$$\ln p(z|a, b) = a \ln b + (-a - 1) \ln z - \ln(\Gamma(a)) - \frac{b}{z}$$

Since the question allows us to omit the terms, which are not depend on z, it could be derived as:

$$\ln p(z|a,b) = (-a-1)\ln z - \frac{b}{z}$$

#### 2.3 Problem 3.2.3

As we know the Bayes theorem:  $p(\theta, \sigma_n^2|y)$  should be,

$$p(\theta, \sigma_{\eta}^{2}|y) = \frac{p(y|\sigma_{\eta}^{2}, \theta)p(\sigma_{\eta}^{2}, \theta)}{p(y)}$$

As  $\sigma_{\eta}^2$  stands for the variance of observational noise, while  $\theta$  represents model parameters, and here, we can assume that they are independent. Hence,

$$p(\sigma_{\eta}^2,\theta) = p(\sigma_{\eta}^2)\dot{p}(\theta)$$

So it will have this formula for posterior,

$$p(\theta, \sigma_{\eta}^{2}|y) = \frac{p(y|\sigma_{\eta}^{2}, \theta)p(\sigma_{\eta}^{2})p(\theta)}{p(y)}$$

Then, taking the logarithm of both sides, we get:

$$\ln p(\theta, \sigma_{\eta}^2 | y) = \ln(p(y | \sigma_{\eta}^2, \theta) p(\sigma_{\eta}^2) p(\theta)) - \ln p(y)$$

Given the distributions:

- Likelihood: $p(y|\sigma_{\eta}^2, \theta)$
- Prior on  $\sigma_{\eta}^2$ :  $p(\sigma_{\eta}^2)$
- Prior on  $\theta$ :  $p(\theta)$

It can be substituted these into the equation:

$$\ln p(\theta, \sigma_{\eta}^2 | y) = \ln p(y | \sigma_{\eta}^2, \theta) + \ln p(\sigma_{\eta}^2) + \ln p(\theta) - \ln p(y)$$

Or even,

$$\ln p(\theta, \sigma_{\eta}^2 | y) \propto \ln p(y | \sigma_{\eta}^2, \theta) + \ln p(\sigma_{\eta}^2) + \ln p(\theta)$$

As p(y) is constant and the goal is to minimize  $lnp(\theta, \sigma_{\theta}^2|y)$ .

 $lnp(sigma_{\eta}^2|a,b)=-(a+1)ln\theta_{\eta}^2-\frac{b}{\theta_{\eta}^2}$  with the previous question. Use what we derived in ex. week 9:

$$\begin{split} lnp(y|\sigma_{\eta}^{2},\theta) &= -\frac{N}{2}ln(2\pi) - \frac{N}{2}ln\sigma_{\eta}^{2} - \frac{1}{2\sigma_{\eta}^{2}}||y - \theta^{T}x||lnp(\theta) \\ &= -\frac{K}{2}ln(2\pi) - \frac{K}{2}ln\sigma_{\theta}^{2} - \frac{1}{2\sigma_{\theta}^{2}}||\theta||^{2} \quad , \ with \ \theta \in R^{k} \end{split}$$

Then,

$$\begin{split} lnp(y|\sigma_{\eta}^{2},\theta) \\ &= -(\frac{K}{2} + \frac{N}{2})ln(2\pi) - \frac{N}{2}ln(\sigma_{\eta}^{2}) - \frac{K}{2}ln(\sigma_{\theta}^{2}) - \frac{1}{2\sigma_{\theta}^{2}}||y - \theta^{T}x|| - \frac{K}{2}ln(\sigma_{\eta}^{2})||\theta||^{2} \\ &\qquad - (a+1)ln(\sigma_{\eta}^{2}) - \frac{b}{\sigma_{\eta}^{2}} \end{split}$$

#### 3 Problem 3.3 Estimation of ICA solution

#### 3.1 Problem 3.3.1

The findings are presented in Figure 5. The computation of the error involves the normalization of columns in both matrices, A and  $\hat{A}$ , followed by potential column switches in  $\hat{A}$  if deemed necessary. Subsequently, the sum of coefficients resulting from the absolute differences between A and  $\hat{A}$  is calculated. An observation of note is the relatively low error, indicating a successful recovery of the sources.

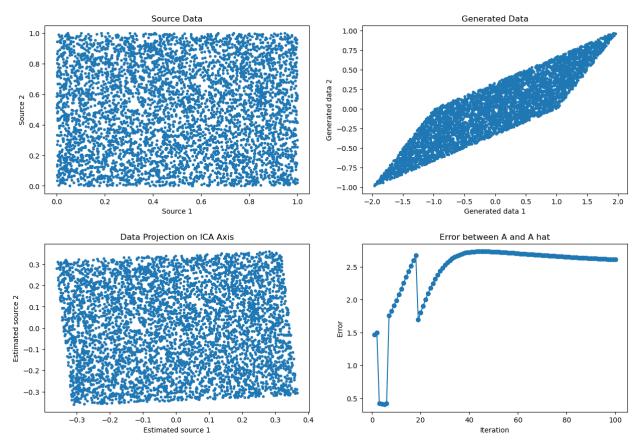


Figure 5: s is drawn from a uniform distribution U(0,1).

## 3.2 Problem 3.3.2

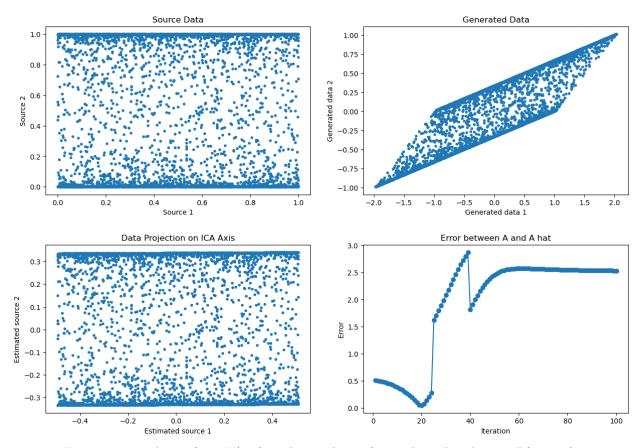


Figure 7: s1 is drawn from U(0;1), and s2 is drawn from a beta distribution B(0.1,0.1).

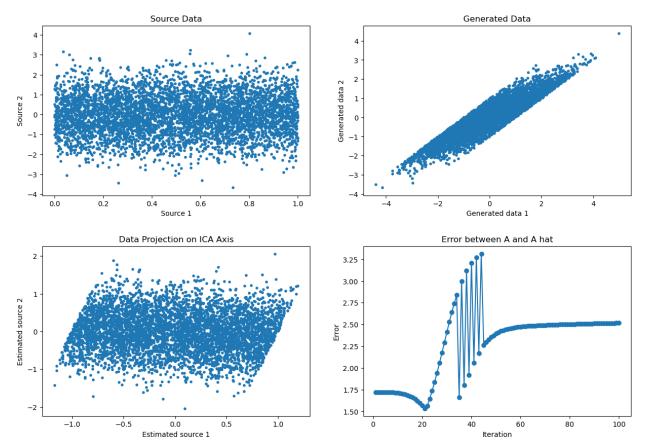


Figure 9: s1 is drawn from U(0,1), and s2 is drawn from a normal distribution N(0,1).

The results are displayed. We can see that ICA is able to successfully estimate in the first two cases but ICA canât unmix the sources for the last case. Indeed, we know from week 8 that we canât unmix gaussian distributed sources with ICA. This can be shown by calculating the distribution of the observations using a change of variables. This leads to the fact that the observation variables and the source variables have the same distribution under the ICA model. Hence, the ICA model is not able to separate Gaussian sources.

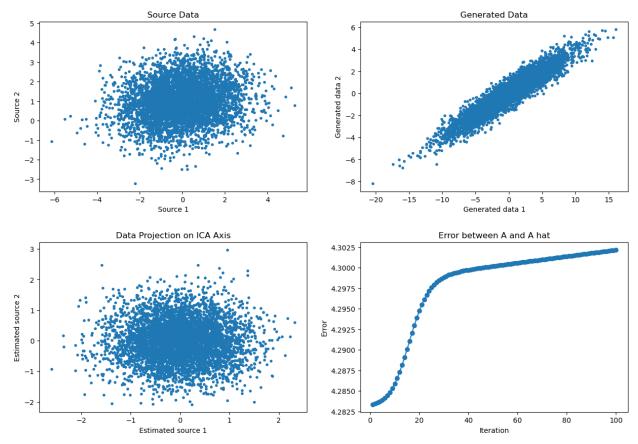


Figure 11: s is drawn from a multivariate normal distribution with  $\mu = (0, 1), \Sigma = [[20.25][0.251]].$ 

#### 4 Problem 3.4 Hidden Markov Models

#### 4.1 Problem 3.4.1

- The number of the states is K=2
- The initial probabilities matrix is  $P_k = \begin{bmatrix} 0.6 \\ 0.4 \end{bmatrix}$
- The transition probabilites matrix is  $P_{ij} = \begin{bmatrix} 0.9 & 0.1 \\ 0.35 & 0.65 \end{bmatrix}$  (As  $P_{11} + P_{12} = P_{21} + P_{22} = 1$ )
- The emission distributions matrix is

$$P(y|k) = \begin{bmatrix} 0.6 & 0.3 & 0.1 \\ 0.1 & 0.6 & 0.3 \end{bmatrix}$$

The sum along each row should be equal to 1.



#### 4.2 Problem 3.4.2

According to ML Sec. 16.5,

$$\alpha(x_1) = P(y_1, x_1) = P(y_1|x_1)P(x_1)$$

As we observe  $a_1$  first:

$$\alpha(x_1) = P(y_1 = a_1|x - 1)P(x_1)$$

Then,

$$\begin{bmatrix} \alpha(x_1 = s_1) \\ \alpha(x_1 = s_2) \end{bmatrix} = \begin{bmatrix} P(y_1 = a_1 | x_1 = s_1) P(x_1 = s_1) \\ P(y_1 = a_1 | x_1 = s_2) P(x_1 = s_2) \end{bmatrix} = \begin{bmatrix} 0.6 \cdot 0.6 \\ 0.1 \cdot 0.4 \end{bmatrix}$$

$$\begin{bmatrix} \alpha(x_1 = s_1) \\ \alpha(x_1 = s_2) \end{bmatrix} = \begin{bmatrix} 0.36 \\ 0.04 \end{bmatrix}$$

As we know,

$$\alpha(x_2) = P(y_2|x_2) \sum_{x_1} \alpha(x_1) P(x_2|x_1)$$

$$= P(y_2|x_2) [\alpha(x_1 = s_1) P(x_2|x_1 = s_1) + \alpha(x_1 = s_2) P(x_2|x_1 = s_2)]$$

Turn into  $a_2$ ,

$$\begin{bmatrix} \alpha(x_2 = s_1) \\ \alpha(x_2 = s_2) \end{bmatrix} = [P(y_2|x_2)[\alpha(x_1 = s_1)P(x_2|x_1 = s_1) + \alpha(x_1 = s_2)P(x_2|x_1 = s_2)]]$$

As we then observe  $a_2$ :

$$\begin{bmatrix} \alpha(x_2 = s_1) \\ \alpha(x_2 = s_2) \end{bmatrix} = \begin{bmatrix} P(y_2 = a_2 | x_2 = s_1) [\alpha(x_1 = s_1) P(x_2 = s_1 | x_1 = s_1) + \alpha(x_1 = s_2) P(x_2 = s_1 | x_1 = s_2)] \\ P(y_2 = a_2 | x_2 = s_2) [\alpha(x_1 = s_1) P(x_2 = s_2 | x_1 = s_1) + \alpha(x_1 = s_2) P(x_2 = s_2 | x_1 = s_2)] \end{bmatrix}$$

$$\begin{bmatrix} \alpha(x_2 = s_1) \\ \alpha(x_2 = s_2) \end{bmatrix} = \begin{bmatrix} 0.3 \cdot [0.36 \cdot 0.9 + 0.04 \cdot 0.35] \\ 0.6 \cdot [0.36 \cdot 0.1 + 0.04 \cdot 0.65] \end{bmatrix} = \begin{bmatrix} 0.101 \\ 0.037 \end{bmatrix}$$

As we know in the ML,

$$P(x_n|y_{[1:n]}) = \frac{\alpha(x_n)}{P(y_{[1:n]})}$$

With 
$$P(y_{[1:n]}) = \sum_{x_n} p(y_{[1:n],x_n}) = \sum_{x_n} \alpha(x_n)$$
 Thus,

$$P(x_2 = s_1 | y_{1:2}) = 0.732$$

#### 4.3 Problem 3.4.3

$$\begin{split} P(y_3|x_2) &= \Sigma_{x_3} P(y_3,x_3|x_2) \quad with \ sum \ rule. \\ &= \Sigma_{x_3} P(y_3|x_3,x_2) P(x_3|x_2) \quad with \ product \ rule \\ &= \Sigma_{x_3} P(y_3|x_3) P(x_3|x_2) \quad because \ of \ the \ graphical \ model \end{split}$$

$$P(y_3 = a_1 | x_2 = s_1) = p(y_3 = a_1 | x_3 = s_1)p(x_3 = s_1 | x_2 = s_1) + p(y_3 = a_1 | x_3 = s_2)p(x_3 = s_2 | x_2 = s_1)$$

$$= 0.6 \cdot 0.9 + 0.1 \cdot 0.1$$

$$= 0.55$$

$$P(y_3 = a_1 | x_2 = s_2) = 0.6 \cdot 0.35 + 0.1 \cdot 0.65 = 0.275$$

#### 4.4 Problem 3.4.4

From ML eq.(16.49) we have:  $P(x_n|y) = \frac{\alpha(x_n)\beta(x_n)}{p(y)}$  and from ML eq.(16.47)  $\beta(x_2) = P(y_3|x_2)$ Thus we get:

$$P(x_2 = s_1 | y_{1:3}) = \frac{\alpha(x_2 = s_1)\beta(x_2 = s_1)}{P(y_{1:3})}$$
$$= \frac{\alpha(x_2 = s_1)\beta(x_2 = s_1)}{\Sigma_{x_3}\alpha(x_3)}$$

$$\beta(x_2 = s_1) = p(y_3 = a_1 | x_2 = s_1) = 0.55$$

As the third observation with Q3 is  $a_1$  Then  $\alpha(x_3) = p(y_3|x_3) \sum_{x_2} p(x_3|x_2) \alpha(x_2)$ 

$$\begin{bmatrix}
\alpha(x_3 = s_1) \\
\alpha(x_3 = s_2)
\end{bmatrix} = \begin{bmatrix}
p(y_3 = a_1 | x_3 = s_1) [p(x_3 = s_1 | x_2 = s_1) \alpha(x_2 = s_1) + p(x_3 = s_1 | x_2 = s_2) \alpha(x_2 = s_2)] \\
p(y_3 = a_1 | x_3 = s_2) [p(x_3 = s_2 | x_2 = s_1) \alpha(x_2 = s_1) + p(x_3 = s_2 | x_2 = s_2) \alpha(x_2 = s_2)]
\end{bmatrix} = \begin{bmatrix}
0.062 \\
0.003
\end{bmatrix}$$

$$\Sigma_{x_3}\alpha(x_3) = \alpha(x_3 = s_1) + \alpha(x_3 = s_2) = 0.062 + 0.03 = 0.065$$
  
Thus  $P(x_2 = s_1|y_{1:3}) = \frac{0.101 \cdot 0.55}{0.065} = 0.855$ 



#### 5 Problem 3.5 Kalman Filter

According to the question ask, the state vector:

$$x_n = \begin{bmatrix} p_n \\ v_n \end{bmatrix}$$

which the  $p_n$  is the position of the time n and the  $v_n$  is the velocity of the object at time n. Here, I use

$$F_n = \begin{bmatrix} 1 & dt \\ 0 & 1 \end{bmatrix}$$

which indicates that the velocity is constant without noise. Besides, I also use dt = 0.1. As the observation is only in the position,  $H_n = \begin{bmatrix} 1 & 0 \end{bmatrix}$ . About  $\eta_n$  and  $v_n$ , I select them such as white noise with a standard deviation of 0.21. The result of the estimation on simulated data is displayed figure 13. The Kalman Filter seems to be able to predict approximately the position of the moving object correctly in the condition.

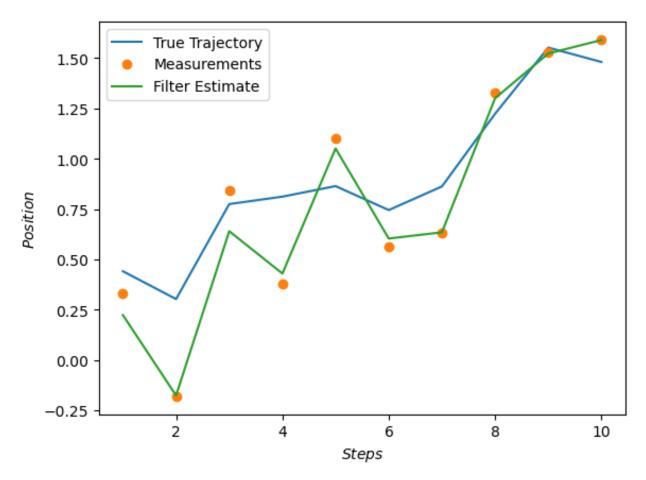


Figure 13: Performance of Kalman Filter for estimating the position of a moving object.

## 6 Problem 3.6 Kernel methods

### 6.1 Problem 3.6.1

We can see in figure 11 and figure 13 that the center location of the chirp is approximately t=5.

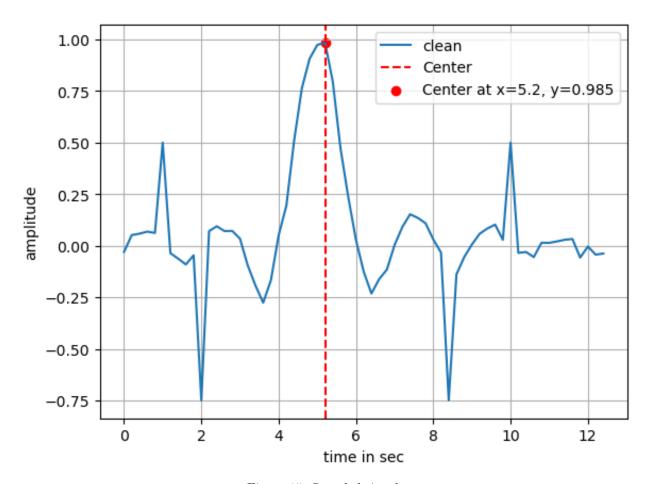


Figure 15: Sampled signal.

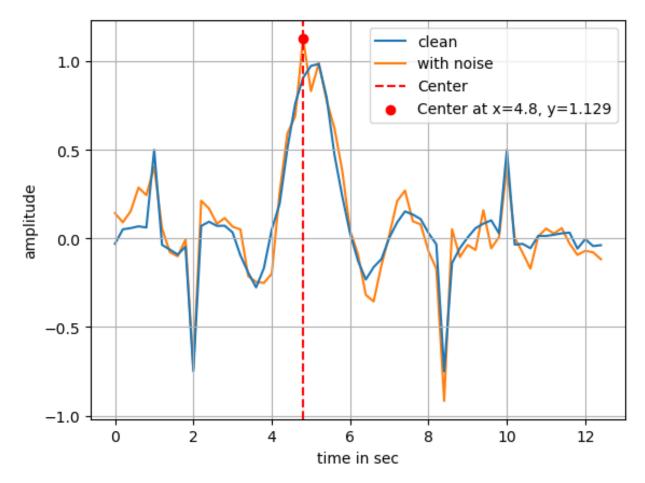


Figure 17: Sampled signal with random white noise.

#### 6.2 Problem 3.6.2

The result of the estimation using Kernel Ridge Regression is available figure 19 and figure 21. I tried different couples of parameters but we can see that the estimation is still not perfect. However, it allows us to get a quite accurate estimation of the center location of the chirp which is approximately t=5. Then, I calculated SNR using the estimated signal and white noise with a standard deviation of 0.004. Indeed, by looking at the data, the noise seems to have a standard deviation smaller than 0.004. Finally, I get SNR = 34.07.

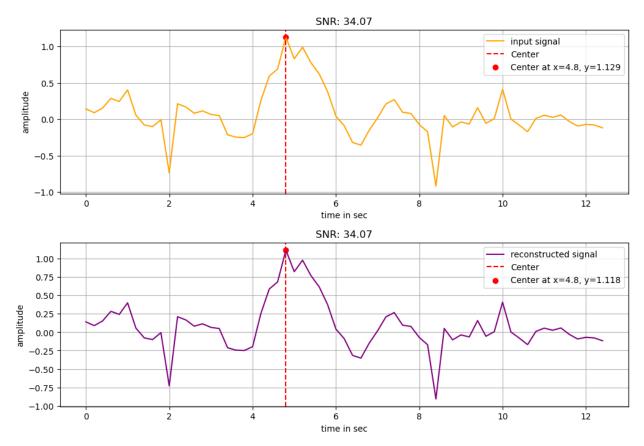


Figure 19: Reconstruct signal using Kernel Ridge Regression. Parameters :  $\sigma=4e-3, C=1e-2.$ 

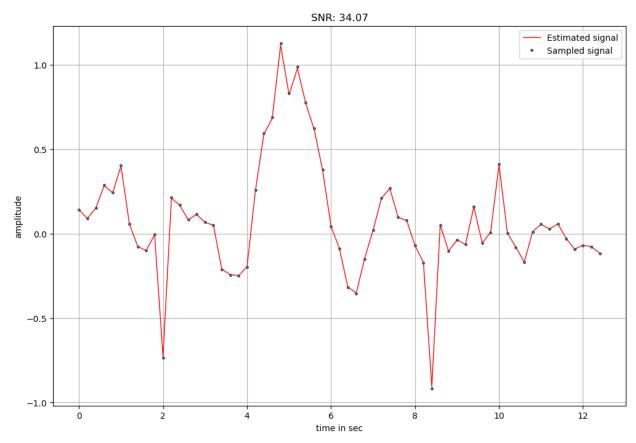


Figure 21: Estimated signal using Kernel Ridge Regression. Parameters :  $\sigma=4e-3, C=1e-2.$ 

#### 6.3 Problem 3.6.3

The result of the estimation using SVR is available figure 23. We can see that SVR seems to perform better than Kernel Ridge Regression. It also allows us to get a quite accurate estimation of the center location of the chirp which is approximately t=5. Then, I calculated SNR using the estimated signal and white noise with a standard deviation of 0.1. Indeed, by looking at the data, I decided that the standard deviation of the noise is actually smaller than 0.1. I get SNR = 47.3. For detecting the outliers, I set all points that are outside of a tube (around the estimate) of a radius 10 times bigger than  $\epsilon$  as outsiders. However, the effect seems not very work on it.

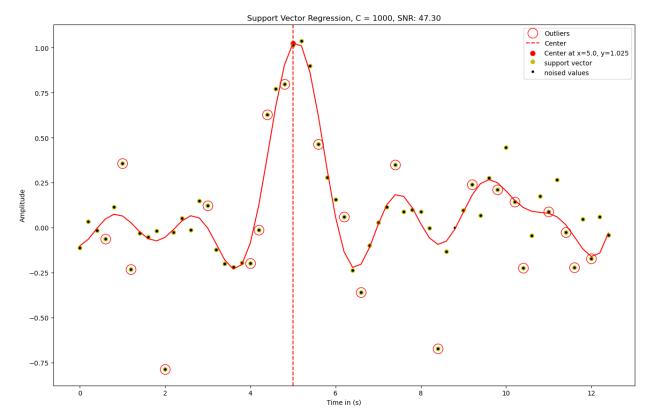


Figure 23: Estimated signal using support vector regression (SVR). Parameters :  $\epsilon=0.01,~\sigma_{kernel}=2,~C=1e3$ 

#### 7 Problem 3.6 Kernel methods

#### 8 code section

#### 8.1 3.1.2

```
Problem_3_1_2.py > ...
       Run Cell | Run Below | Debug Cell
       #%%
  1
  2
       import numpy as np
       import matplotlib.pyplot as plt
       from sklearn.linear_model import Lasso
  4
       from scipy.fft import dct, idct
  6
       from scipy.io import loadmat
       Run Cell | Run Above | Debug Cell
  8
       #%%
       #Load data from problem3_1.mat
       data = loadmat('/Users/luchengliang/ML_sp/Problem Set 3/probl
 10
 11
       n = data['n'].flatten()
       x = data['x'].flatten()
 12
 13
       Run Cell | Run Above | Debug Cell
 14
       #%%
 15
       # Signal setup
       N = 2**5 # number of observations to make
 16
 17
       l = 2**9 # signal length
 18
 19
       B = np.zeros((N, l))
       for i in range(N):
 20
 21
           B[i, n[i]] = 1
 22
 23
       # Since it is sparse in the IDCT domain, i.e. B*x = B*Phi*X
       # where X sparse, BF = B*Phi; and Phi is the DCT matrix, Phi
 24
 25
       # Equivalently, since IDCT = transpose of DCT using idct we d
       BF = idct(B, norm='ortho')
 26
 27
 28
       lambda_ = 0.005
       model = Lasso(lambda_, fit_intercept=False) University of Denmark
 29
       model.fit(BF, x)
 30
       solsB = model.coef_
 31
 32
 33
       # create IST solution
       nstens = 100000
 34
```

```
t_{:, k} = np.sign(t_{tilde})*np.maximum(abs(t_{tilde}) - lambda_*mu, 0)
     solsIST = t_{:, -1}
     sols = solsIST*np.sqrt(2 / len(solsIST)) # normalize according to DCT specification
     a = sols[np.nonzero(sols)] # get ai values
45
     k = np.count_nonzero(sols) # get number of non-zero components in DCT domain
     m = np.where(sols)[0] # get positions of non-zero components
     print("a_i values:", a)
     print("K status:", k)
     print("m_j values:", m)
     # Reconstruct the multitone signal for verification
     t = np.arange(0, l) # time axis
     s_multitone = np.zeros(l)
     for i in range(k):
         s_{multitone} += a[i] * np.cos((np.pi * (2 * m[i] - 1) * t) / (2 * l))
     # `s_multitone` now contains the reconstructed multitone signal
     # plot solutions
     fig, ax= plt.subplots(1, 1, figsize=(6, 3))
     ax.stem(solsB, markerfmt='bo', label='sklearn Lasso', basefmt=' ')
     ax.stem(solsIST, markerfmt='ro', label='IST', basefmt=' ')
     ax.legend()
     ax.set_title('Solutions')
     # Take the inverse IDCT (i.e. the DCT) in order to compute the estimated signal.
     x_hat = dct(solsIST, norm='ortho')
     b_hat = dct(solsB, norm='ortho')
     fig, ax= plt.subplots(3, 1, figsize=(12, 8))
     ax[0].plot(b_hat)
     ax[0].plot(n, x, 'r.')
     ax[0].set_title('Estimated in time domain by LASSO')
     ax[1].plot(x_hat)
     ax[1].plot(n, x, 'r.')
     ax[1].set_title('Estimated in time domain by IST')
     ax[2].plot(s_multitone)
     ax[2].set_title('Estimated signal reconstructed using multitone formula')
     fig.tight_layout()
     plt.show()
```

#### 8.2 3.3

```
#%%
     import numpy as np
     import matplotlib.pyplot as plt
     from sklearn.decomposition import PCA
     Run Cell | Run Above | Debug Cell
     #%%
     def column_switch(A, A_hat):
          # Compute differences and normalize columns
          diff = np.abs(A - A_hat)
          norm_diff = diff / np.linalg.norm(diff, axis=0)
          # Check if columns need to be switched in A_hat
          for i in range(A.shape[1] - 1):
              if np.sum(norm_diff[:, i]) > np.sum(norm_diff[:, i + 1]):
                  # Switch columns in A_hat
                  A_hat[:, [i, i + 1]] = A_hat[:, [i + 1, i]]
          return A_hat
     Run Cell | Run Above | Debug Cell
     def error_cal(A, A_hat):
          #Check if columns need to be switched in A_hat
24
          A_hat = column_switch(A, A_hat)
          # Normalize columns of A and A_hat
          A_normalized = A / np.linalg.norm(A, axis=0)
          A_hat_normalized = A_hat / np.linalg.norm(A_hat, axis=0)
30
          # Compute the absolute difference between normalized matrices
          diff = np.abs(A_normalized - A_hat_normalized)
          # Compute the sum of coefficients from the absolute difference
          error = np.sum(diff)
          return error
     Run Cell | Run Above | Debug Cell
     # Independent Component Analysis function
     def ICA(x, mu, num_components, iters, mode, A):
```

```
# Random initialization
          W = np.random.rand(num_components, num_components)
         N = np.size(x, 0)
          if mode=='superGauss':
              phi = lambda u : 2*np.tanh(u)
          elif mode=='subGauss':
47
              phi = lambda u : u-np.tanh(u)
          else:
              print("Unknown mode")
              return W
          errors = []
          for i in range(iters):
              u = W @ x.T
              dW = (np.eye(num_components) - phi(u) @ u.T/N) @ W
              # Update
              W = W + mu*dW
              A_hat = W.T
              error = error_cal(A, A_hat)
              errors.append(error)
          return W, errors
     Run Cell | Run Above | Debug Cell
     def Find_the_Source(s):
          # Mix signals
70
          A = np.array([[3, 1], [1, 1]])
          x = (A@s).T
          r = s.T
          # calculate ica
          mu = 0.1
          components = 2
          iterations = 100
79
          # Mean across the first (column) axis
          col_means = np.mean(x, axis=0)
```



```
x = x - col_means
           # run ICA
           W, errors = ICA(x, mu, components, iterations, 'subGauss', A)
           # Normalize unmixing matrix
           W = np.divide(W, np.max(W))
           # Compute unmixed signals
           y = (W@x.T).T
           # Plotting
           plt.figure(figsize=(15, 10))
           plt.subplot(2, 2, 1) # Subplot 1: Source Data
           plt.plot(r[:, 0], r[:, 1], '.')
           plt.xlabel('Source 1')
           plt.ylabel('Source 2')
100
           plt.title('Source Data')
           plt.subplot(2, 2, 2) # Subplot 2: Generated Data
           plt.plot(x[:, 0], x[:, 1], '.')
104
           plt.xlabel('Generated data 1')
           plt.ylabel('Generated data 2')
106
           plt.title('Generated Data')
           plt.subplot(2, 2, 3) # Subplot 3: Data Projection on ICA Axis
           plt.plot(y[:, 0], y[:, 1], '.')
           plt.xlabel('Estimated source 1')
110
111
           plt.ylabel('Estimated source 2')
112
           plt.title('Data Projection on ICA Axis')
113
114
           iters = np.arange(1, iterations + 1)
           plt.subplot(2, 2, 4) # Subplot 4: Error by Iteration
116
           plt.plot(iters, errors, marker='o')
           plt.xlabel('Iteration')
118
           plt.ylabel('Error')
119
           plt.title('Error between A and A hat')
120
121
           plt.subplots_adjust(hspace=0.3)
122
          plt.show()
```



```
Run Cell | Run Above | Debug Cell
124
125
      # generate data
126
127
      N = 5000
128
129
      # Define two non-gaussian uniform components
130
      s1 = np.random.rand(N)
      s2 = np.random.rand(N)
131
132
      s = np.array(([s1, s2]))
133
134
      # Define one non-gaussian uniform component and one beta component
135
      s1b = np.random.rand(N)
136
      s2b = np.random.beta(0.1, 0.1, size=N)
137
      sb = np.array(([s1b, s2b]))
138
139
      # Define one non-gaussian uniform component and one gaussian component
140
      s1n = np.random.rand(N)
141
      s2n = np.random.normal(size=N)
142
      sn = np.array(([s1n, s2n]))
143
144
      #Define multivariate normal distribution with
145
      \#\mu = (0, 1), \Sigma = [2 \ 0.25; \ 0.25 \ 1]
146
      mean = [0, 1]
147
      covariance = [[2, 0.25], [0.25, 1]]
148
      sm_r = np.random.multivariate_normal(mean, covariance, N)
149
      sm = sm_r.T
150
151
      Find_the_Source(s)
152
      Find_the_Source(sb)
153
      Find_the_Source(sn)
154
      Find_the_Source(sm)
155
```

#### 8.3 3.5

```
1
     import numpy as np
     import matplotlib.pyplot as plt
     # Set the parameters
     q = 1
     dt = 0.1
     s = 0.21
10
     F = np.array([
         [1, dt],
11
         [0, 1],
12
13
     1)
14
     Q = q*np.array([
15
          [s**2, 0],
          [0, 0]
17
     1)
19
     H = np.array([1, 0])
20
     R = s**2*np.identity(2)
21
     m0 = np.array([[0], [1]])
22
     P0 = np.identity(2)
23
```

```
24
     # Simulate data
25
26
     np.random.seed(1)
27
28
     steps = 10
29
     X = np.zeros((len(F), steps))
     Y = np.zeros((len(H), steps))
30
     x = m0
31
     for k in range(steps):
32
         x = F@x + s*np.random.randn(len(F), 1)
33
         y = H@x + s*np.random.randn(1, 1)
34
         X[:, k] = x[:, 0]
35
         Y[:, k] = y[:, 0]
36
37
38
39
     # Kalman filter
40
41
     m = m0
```

```
P = P0
42
     kf_m = np.zeros((120/m) - V.chopo[1]))
43
     kf_P = np.zeros(( (property) shape: _Shape ; [1]))
44
     for k in range(Y.shape[1]):
         m = F@m
47
         P = F@P@F.T + Q
         e = Y[:, k].reshape(-1, 1) - H@m
         S = H@P@H.T + R
50
51
         K = P@H.T@np.linalg.inv(S)
52
         m = m + K@e
53
         P = P - K@S@K.T
54
         kf_m[:, k] = m[:, 0]
         kf_P[:, :, k] = P
56
57
58
     a = np.arange(1,11)
59
     plt.figure()
     plt.plot(a, X[0, :], '-')
60
     plt.plot(a, Y[0, :], 'o')
61
     plt.plot(a, kf_m[0, :], '-')
62
     plt.legend(['True Trajectory', 'Measurements', 'Filter Estimate'])
63
     plt.xlabel('$Steps$')
64
     plt.ylabel('$Position$')
66
67
     plt.figure()
68
     plt.plot(X[0, :], X[1, :], '-')
     plt.plot(Y[0, :], Y[1, :], 'o')
69
70
     plt.plot(kf_m[0, :], kf_m[1, :], '-')
     plt.legend(['True Trajectory', 'Measurements', 'Filter Estimate'])
71
72
     plt.xlabel('$x_1$')
     plt.ylabel('$x_2$')
73
74
```



#### 8.4 3.6

```
#%%
    Run Cell | Run Above | Debug Cell
2 ∨ #%%
3 ∨ import os
    from scipy.io import loadmat
    import numpy as np
    import soundfile as sf
    import matplotlib.pyplot as plt
    Run Cell | Run Above | Debug Cell
8 \( \square\) #\%
    data = loadmat('./problem3_6.mat')
    t = data['t'].flatten()
    y = data['y'].flatten()
    # Calculate the center of the signal
    max_value = np.max(y)
    ind = np.argmax(y)
    tmax = t[ind]
    max_value_abbrev = round(max_value, 3)
    #Initial figure and plot the signal and its center point
    fig, ax = plt.subplots()
    ax.plot(t, y, label='clean')
    ax.axvline(tmax, color='r', linestyle='--', label='Center')
    ax.scatter(tmax, max_value, color='red', label=f'Center at x={tmax}, y={max_value_abbrev}')
    ax.set_xlabel('time in sec')
    ax.set_ylabel('amplitude')
    ax.legend()
    ax.grid()
    #Initial another figure
    fig, ax = plt.subplots()
    y_original = y
    ax.plot(t, y_original, label='clean', zorder=1)
    np.random.seed(0)
    N = t.size
```

```
percent_outlier = 0.1
sigma = 0.004
noise = np.random.randn(N)
noise *= (np.sum(y**2)/np.sum(noise**2)/10**(snr/10))**0.5
y += noise
ax.plot(t, y, label='with noise', zorder=0)
max_value_mednoise = np.max(y)
ind_mednoise = np.argmax(y)
tmax_mednoise = t[ind_mednoise]
tmax_mednoise_abbrev = round(tmax_mednoise, 1)
max_value_mednoise_abbrev = round(max_value_mednoise, 3)
ax.axvline(tmax_mednoise, color='r', linestyle='--', label='Center')
ax.scatter(tmax_mednoise, max_value_mednoise, color='red', label=f'Center at x={tmax_mednoise_abbrev}, y={max_value_mednoise_abbrev}')
ax.set_xlabel('time in sec')
ax.set_ylabel('amplitude')
ax.legend()
ax.grid()
Run Cell | Run Above | Debug Cell
pair_dist = np.abs(t.reshape(-1, 1) - t.reshape(1, -1))
K = np.exp(-1/(sigma**2)*pair_dist**2)
A = C*np.identity(N) + K
samples = t[-1]
x = np.arange(0, samples + 0.2, 0.2)
M2 = len(x)
z0 = np.zeros(M2)
     z0[k] = 0
     for j in range(N):
    value = np.exp(-1/(sigma**2)*(t[j] - x[k])**2)
         z0[k] += sol[j]*value
```



```
max_value_re = np.max(z0)
ind_re = np.argmax(z0)
 tmax_re = x[ind_re]
 tmax_re_abbrev = round(tmax_re, 1)
max_value_re_abbrev = round(max_value_re, 3)
SNR = np.var(z0) / (0.2 ** 2 * np.var(np.random.rand(1, len(z0))))
 fig, (ax1, ax2) = plt.subplots(2, 1, figsize=(12, 8))
 ax1.set_xlabel('time in sec')
ax1.set_ylabel('amplitude')
ax1.plot(t, y, label='input signal', color='orange')
ax1.axvline(tmax_mednoise, color='r', linestyle='--', label='Center')
ax1.scatter(tmax_mednoise, max_value_mednoise, color='red', label=f'Center at x={tmax_mednoise_abbrev}, y={max_value_mednoise_abbrev}')
ax1.legend()
ax1.set_title(f'SNR: {SNR:.2f}')
 ax1.grid()
ax2.set_xlabel('time in sec')
ax2.set_vlabel('amplitude')
ax2.plot(x, z0, label='reconstructed signal', color='purple')
ax2.axvline(tmax_re, color='r', linestyle='--', label='Center')
ax2.scatter(tmax_re, max_value_re, color='red', label=f'Center at x={tmax_re_abbrev}, y={max_value_re_abbrev}')
ax2.set_title(f'SNR: {SNR:.2f}')
ax2.grid()
plt.subplots_adjust(hspace=0.3)
 fig, ax = plt.subplots(figsize=(12, 8))
ax.set_xlabel('time in sec')
ax.set_ylabel('amplitude')
ax.plot(x, z0, 'r', linewidth=1, label='Estimated signal')
ax.plot(t, y, '.', markeredgecolor=0.3*np.array([1, 1, 1]), markersize=5, label='Sampled signal')
 ax.legend()
 ax.set_title(f'SNR: {SNR:.2f}')
 ax.grid()
```



```
import numpy as np
     from scipy.io import loadmat
     import librosa
     import librosa.display
      🔓 port matplotlib.pyplot as plt
     from random import randrange
     from sklearn.svm import SVR
     Run Cell | Run Above | Debug Cell
     #%%
10
     def awgn(signal, snr):
         x_watts = signal ** 2
         # Set a target SNR
         target_snr_db = snr
13
         # Calculate signal power and convert to dB
         sig_avg_watts = np.mean(x_watts)
         sig_avg_db = 10 * np.log10(sig_avg_watts)
         noise_avg_db = sig_avg_db - target_snr_db
         noise_avg_watts = 10 ** (noise_avg_db / 10)
         # Generate an sample of white noise
         mean noise = 0
         noise_volts = np.random.normal(mean_noise, np.sqrt(noise_avg_watts), len(x_watts))
23
         # Noise up the original signal
         y_volts = signal + noise_volts
         return y_volts
     #Load the data
     data = loadmat('./problem3_6.mat')
     t = data['t'].flatten()
     y = data['y'].flatten()
     x = t
     # parameters
     N=t.size
     snr = 10 \#dB
     percent_outlier = 0.1
38
     # learning parameters
     epsilon=0.01
     kernel_type='Gaussian'
```



```
kernel_params=2
      y_noised = awgn(y, snr)
      x_{col} = x.reshape((np.size(x), 1))
     y_row = np.copy(y_noised)
t_col = t.reshape(( np.size(t), 1))
     t_col = np.around(t_col, decimals=4)
x_col = np.around(x_col, decimals=4)
     y_row = np.around(y_row, decimals=4)
     gamma = 1/(np.square(kernel_params)) # gamma needs to be calculated in order to use 'Gaussian' kernel, which is not available in the library regressor = SVR(kernel='rbf', gamma=gamma, C=C, epsilon=epsilon)
      regressor.fit(x_col,y_row)
61
62
      y_pred = regressor.predict(t_col)
     threshold = 10 * epsilon
65
66
      outsider = np.zeros(len(t))
      for i, sv_index in enumerate(regressor.support_):
        j = np.where(x_col == x_col[sv_index])[0][0]
         if abs(y_row[sv_index] - y_pred[j]) > threshold:
    outsider[i] = 1
     outsider = outsider.astype(bool)
     max_value = np.max(y_pred)
      ind = np.argmax(y_pred)
     t_{max} = x[ind]
     tmax_abbrev = round(t_max, 1)
      max_value_abbrev = round(max_value, 3)
      SNR = np.var(y_pred) / np.var(0.2 ** 2 * np.random.randn(len(y_pred)))
```

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