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Deep Clustering for ANN Supported Source Separation

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1 Introduction

When perceiving sound created by multiple acoustic sources, the human auditory system is extremely good at focussing on parts of the sound coming from only one acoustic source.

Imagine yourself at a cocktail party. The sound that you perceive is made up of all kinds of acoustic sources: The voice of the person you are listening to, the sound of music, the background chatter of other people talking, the sound of clinking glasses, etc. Still you can clearly understand the speaker in your group, even though its acoustic signal overlaps at every moment with the acoustic signal of other people speaking and background noise.

The difficulty of understanding speech in a multiple speaker environment is often called "The cocktail party problem" [4], first mentioned by Colin Cherry in 1953. It was also Colin Cherry, who first brought up the idea of automatic speech separation in his famous paper on "experiments of the recognition of speech" [7]. The first method to yield some promising results was introduced by A.S. Bregman in 1990 using *Auditory scene analysis* [3]. Since then, multiple methods have been explored. One of them is *Spectral Clustering*, a method based on partitioning data points into clusters according to eigenstructure of a similarity matrix.

With the recent rise of deep learning in a variety of applications, the first method to be applied to source separation was presented by Chao Weng and co. in 2015 [20]. Quickly, different deep learning methods based on a technique called *Deep Clustering* emerged, forming the state-of-art in the source separation problem. This report, will focus on the deep clustering method as it was introduced by John R. Hershey and co. in 2015 [10] and will study *TasNet*, a deep clustering system performing audio separation in the time domain yielding state-of-the art results [12].

The applications of automated source separation are wide-ranging. To name a few:

- *Automatic meeting transcription*: In business meetings, multiple speakers are alternating in a short time or even speaking at the same time. Making it possible to automatically separate speakers from each other and transcribe would be of great use for companies.
- *Virtual assistant*: Virtual assistants, like Alexa, Siri and Google Home are playing a critical part in smart home systems. Filtering out noise and separating the owner's voice from other voices are challenges that can be taken on by automated source separation techniques.
- *Automatic subtitling of music/video*: According to the World Health Organization (WHO) over 5% of the world population suffers from disabling hearing loss [1]. Subtitling of music and video content, which are often made up of sound coming from multiple acoustic sources, is therefore essential for them.

To begin with, the source separation problem is mathematically defined and basic mathematical operations and basic deep learning methods used in deep clustering are explained 2. In the following conventional methods and first attempts of using deep learning techniques for automated source separation are presented 3. The essence of this paper being deep clustering and *TasNet* will be presented in-detail in section 4. Finally, a conclusion is drawn 5.

2 Basics

This section firstly defines the source separation problem and then gives an overview of the basic operations and methods used in systems.

2.1 Definition source separation problem

A general definition of the source separation problem was given by Jean-Francois Cardoso, defining the problem as “recovering unobserved signals or sources from several observed mixtures” [5]. He also added two important restrictions that we will apply to our definition as well:

1. the source signals are not observed
2. no information is available about the mixture

The condition of not knowing how many sources the mixture is composed of, became known as *blind* source separation [5]. In this paper, we add another restriction: The signal is observed over a single channel (single microphone). To sum it up, the *source separation problem* is defined in this paper as “recovering the original signals of an unknown amount of sources from a single-channel mixture of the source signals”.

Putting the above definition in a mathematical form:

- the unknown number of acoustic sources: N .
- the original sound of the i -th source at time t : $x_i(t)$.
- the single received sound signal at time t : $X(t) = \sum_{i=1}^N x_i(t)$.

The goal of source separation is to recover $x_1(t), \dots, x_N(t)$ from $X(t)$. An algorithm performing this task is presented below.

Input: $X(t)$

Output: $(N, \{x_1(t), \dots, x_N(t)\})$

2.2 Mathematical operations

This subsection aims to shine light on essential mathematical operations used in the systems that will be explained in sections 3, 4.

2.2.1 Short term fourier transform

In order to understand the short term fourier transform, one firstly has to get an understanding of the fourier transform itself. Having a continuous signal $s(t)$, the fourier transform is defined as

$$S(f) = \int_{\mathbb{R}} s(t) e^{-2j\pi ft} dt$$

The fourier transform $S(f)$ gives the magnitude of every frequency in the signal. A signal can be decomposed as the sum of signals consisting only of one frequency (so called sinusoids). Thus, one can say that the higher the magnitude of a frequency of the fourier

transform, the higher is the percentage of the signal being composed of the sinusoid of this frequency. An important aspect to notice is that the fourier transform integrates over the whole time domain of the signal in order to get the exact magnitude of every frequency present in the signal. Thus integrating only over a part of the time domain leads to approximated values of the frequencies magnitudes. On the other hand, a shorter time interval to integrate over leads to a better time localization of the frequency. This famous trade-off is well-known as the uncertainty principle in signal analysis [14].

The short term fourier transform divides the signal into local sections using a window function $w(\tau) = 0 : \tau \notin [-a, a]$ over which the fourier transform is applied.

$$S(f, t) = \int_{\mathbb{R}} s(\tau)w(\tau - t)e^{-2j\pi f\tau}d\tau$$

Applying the short term fourier transform converts a signal $X(t)$ to multiple (F, T) bins that can be plotted in a *spectrogram*. It therefore is used to describe the change in frequency over time. Using an appropriate window function leads to a good trade off between time localization and frequency approximation.

2.2.2 Embedding space

An embedding is a structure-preserving, injective mapping $f : X \rightarrow Y$. Therefore, every object of domain X is mapped to a distinct object in domain Y . In this report, we are interested in embeddings that map to normed spaces meaning that it is possible to measure the similarity of two objects x_1 and $x_2 \in X$ by measuring its distance $d(y_1; y_2) \in \mathbb{R} : y_1, y_2 \in Y$. It is essentially used to convert data to a feature representation where certain properties can be preserved by distance, so that the data is easily comparable by computers.

One famous example are word embedding models, such as *word2vec* that maps words to a vector representation. Assuming that words have multiple degrees of similarity, word embeddings are used to easily calculate the similarity of words on multiple axis [13]. Later, we will see that the function mapping objects from one domain to an embedding space can be trained on artificial neural networks and are very powerful tools for the source separation problem 4.1.

2.3 Deep neural network basics

This subsection aims to clarify basic deep learning methods and architectures used in the sections 3, 4.

2.3.1 Convolutional neural network

Convolutional neural networks are neural networks that are specialized for processing data that has grid-like topology, for example time series data for a 1D convolution [9]. The idea is based on the convolution operation $*$ being:

$$s(t) * g(t) = \int_{\mathbb{R}} s(\tau)g(t - \tau)d\tau$$

which corresponds to

$$s(n) * g(n) = \sum_{i=-\infty}^{\infty} s(i)g(n-i)$$

in the discrete case. The fourier transform of a signal to its spectrum can essentially be seen as a convolution operation [21].

This principle can now be applied to artificial neural networks. Let x be the output layer and w the weights of a convolutional layer. Such a layer is defined by its filter size k and other parameters, such as stride, padding, etc. For more information please refer to [9]. The weight array is also called “Kernel” [19] and normally a convolutional layer is decomposed of multiple kernels, such that there are multiple parallel outputs. The convolution operation in neural networks is defined as follows:

$$y_{i,j} = w_i * x = f\left(\sum_{n=j}^{j+k} (w_{i,n-j}x_n) + b_i\right)$$

$y_{i,j}$ presents the j -th neuron of the output layer of the i -th kernel. Similarly $w_{i,l}$ presents the l -th weight of the i -th kernel. x_n is the output of the n -th neuron of the previous layer and b_i is the bias corresponding to the i -th kernel. Finally $f()$ is the activation function applied to each convolution operation.

Convolutional neural networks can very well be applied to act as filterbanks for the signal as was shown in [8] and as we will later see in subsection 4.1.

2.3.2 Recurrent neural network

Recurrent neural networks are neural networks that are “specialized in processing a sequence of values x_1, \dots, x_t ” [9]. For that reason, they are very well applicable when dealing with acoustic signals having high autocorrelation (high correlation between $s(t)$ and $s(t-k, s(t+k))$ for $k > 0$). Recurrent neural networks are extremely effective when dealing with sequential data. First, they process data of different time instances in one computation step, thus taking full advantage of autocorrelation. Second, they can handle data sequences of variable length in contrast to feed forward neural networks where the input length is determined by the size of the input layer. This is achieved by loops between recurrent units inside the neural network [9]. Unfolding a recurrent unit z of a recurrent neural network at different time instances leads to so-called hidden copies of the network being connected to each other: $z^0, z^1, \dots, z^{t-1}, z^t$. As a result, such a recurrent unit z^t at time t takes both the output of the node leading to it (x^t) and the output of the same recurrent unit of one time instance earlier (z^{t-1}) as input resulting in:

$$z^t = f(Vx^t + Wz^{t-1} + b)$$

with $f()$ being the activation function, V the weights between recurrent unit z and input x , W the weights between two hidden recurrent units z^{t-1} and z^t and b the constant bias term. This basic setup is known to have the vanishing/exploding gradient problem due to an exponential decrease/increase in value of long-term components ($\frac{\partial z^t}{\partial z^{t-k}} \rightarrow \text{const.} \times W^k \rightarrow 0$ or ∞) [15]. To overcome this problem, Hochreiter and co. introduced a “better” recurrent unit, called long shot term memory.

2.3.3 Long short term memory

To overcome the exploding/vanishing gradients problem, the long short term memory unit introduces an internal state which can be seen as the unit's "memory" [11]. At every time instance t , the unit updates its memory c^t using the memory c^{t-1} and the unit's output h^{t-1} , as well as the new input x^t . c^t in combination with h^{t-1} and x^t is then used to produce the new output h^t . The structure of the long short term memory unit is shown in Figure 1. It is important to notice that the long short term memory uses x^t and h^{t-1} to compute four different gates: f, i, g, o which are used to control the flow of information in the unit. Also, we can see that when doing backpropagation, the gradient easily "flows" to earlier hidden units without the weight matrix being applied to it (shown by the red arrow) in 1. This "highway" for the gradient strongly reduces the vanishing/exploding gradient effect and allows the unit to take into account long term dependencies of the input data.

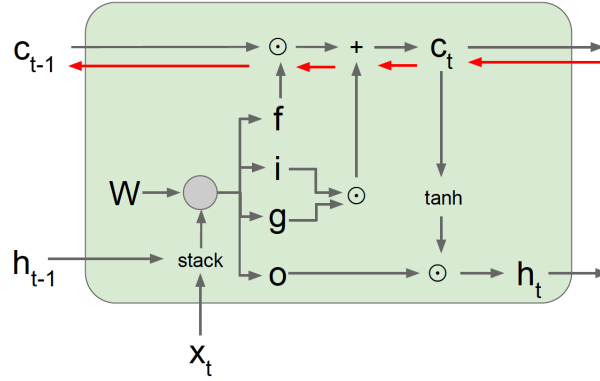


Figure 1: Long short term memory

To further improve this ability, bi-directional long short term memory units were proposed to exploit correlations between data in both time directions and are used in many state-of-the-art systems nowadays. Going more into detail is out of scope for this report, but more in-detail information can be found in [11].

2.3.4 Autoencoder

An autoencoder is a neural network that is "trained to attempt to copy its input to its output" [1]. It always consists of two parts, the encoder and the decoder. In its simplest form, the encoder maps the input x to one hidden layer $f(x) = z$ and the decoder tries to reconstruct the original data $g(z) = x'$ from this layer. Even though autoencoders are trained with loss functions trying to minimize the difference between x and x' , perfect reconstruction of the data input is not the goal. Instead, autoencoder should be trained in a way preventing them to just copy the input to the output, but to extract useful information about data distribution instead [9]. Thereby, the hidden layer should represent the most useful information. Two approaches are considered in this report. First, an undercomplete autoencoder is defined by the hidden layer having much lower dimension than the data, thus forcing the network to "store all data in a compressed form". A drawback of this approach is that the hidden layer might be too small to learn enough useful information. Second, a sparse autoencoder forces the hidden layer to be sparse, by

introducing a sparsity penalty on the hidden layer $P(h)$, which will then be added to the loss function $\mathbb{L}(x, x') + P(h)$. Especially in the field of machine translation, autoencoders in combination with recurrent neural networks produce state-of-the-art results [18]. For more detailed information, please refer to [9].

3 Conventional methods

The first theoretical framework for automated source separation was introduced by A.S. Bregman in 1990 [3]. Systems using this approach are called *computational auditory scene analysis* and produced first results, that are acceptable [16]. Section 3.1 will handle cover the system.

Spectral Clustering, a method well-know for data clustering, was then applied to the source separation problem and showed good results (see 3.2). It was the first method used for source separation to train some of its parameters on actual data [2].

3.1 Computational auditory scene analysis

Computational auditory scene analysis for source separation is an automation of the auditory scene analysis which is inspired by the principles of processing sound in human auditory system [16]. In computational auditory scene anlysis the algorithm consists of two parts: *segmentation* and *grouping*. A more ex A detailed description of the algorithm can be seen in the following and it based on the paper of *Wang* and *Brown* [16].

Input: $(N, X(t))$

Algorithm:

1. Segmentation:
 - $X(t)$ is transformed into time-frequency space using short term fourier transform
 - Segmenation rules are used to define different time-frequency (T,F) regions.
2. Grouping:
 - (F, T) bins are combined to streams corresponding to sound sources
 - Auditory masks are created using different streams

Output: $(\{x_1(t), \dots, x_N(t)\})$

The segmentation rules used in the segmentation part are hand-crafted similarity features upon which the different (T,F) regions are created. For every speaker, an audiotry masks is created as a last step of the grouping part. These auditory masks are boolean matrices that can be applied to the spectrogram to retrieve the part of the spectrogram corresponding to a speaker. Important to notice is that the amount of different speakers N has to be an input of the algorithm and cannot be computed by the system itself. Since computational auditory scene analysis is a rule based system it is very difficult to generalize to broader conditions and needs very precise tuning in order to work.

Computational auditor scene analysis was the first version to produce some meaningful results, but is not powerful enough for the general source seperation problem as we defined

it in section 2.1.

3.2 Spectral Clustering

Spectral Clustering is a well-known method in multivariate statistic for the clustering of data. It makes use of the eigenvalues (also called spectrum, thus the name *Spectral clustering*) of a similarity matrix W of the data to reduce the dimensionality of the data. Since spectral clustering will be very important in latter chapters it will be described here in detail. The similarity matrix W can be defined in multiple ways depending on the task, but it should have the following properties.

1. W should be a hermitian matrix: $W^T = W$.
2. $(W)_{i,j} \geq 0, \forall i, j$
3. W should be non-negative definite.

Considering a dataset of P entries $D \in \mathbb{R}^P$, describes the similarity between each of these datapoints, thus $W \in \mathbb{R}^{P \times P}$. Since W can be considered as a distance measure, whereas $W(d_1, d_2) = W(d_2, d_1$ and with only positive distances, the first two of the above defined properties seem reasonable. The third property ensures that all eigenvalues of W are ≥ 0 being a characteristic of non-negative definite matrices.

Having defined a similarity matrix, an example of an algorithm for spectral clustering used in [2] is presented in figure 2.

Input: Similarity matrix $W \in \mathbb{R}^{P \times P}$.

Algorithm:

1. Compute first R eigenvectors U of $D^{-1/2}WD^{-1/2}$ where $D = \text{diag}(W\mathbf{1})$.
2. Let $U = (u_1, \dots, u_P)^T \in \mathbb{R}^{P \times R}$ and $d_p = D_{pp}$.
3. Initialize partition A .
4. Weighted K -means: While partition A is not stationary,
 - a. For all r , $\mu_r = \sum_{p \in A_r} d_p^{1/2} u_p / \sum_{p \in A_r} d_p$
 - b. For all p , assign p to A_r where $r = \arg \min_{r'} \|u_p d_p^{-1/2} - \mu_{r'}\|$

Output: partition A , distortion measure $\sum_r \sum_{p \in A_r} d_p \|u_p d_p^{-1/2} - \mu_r\|^2$

Figure 1: Spectral clustering algorithm that minimizes $J_1(W, E)$ with respect to E with weighted K -mean. See Section 2.6 for the initialization of the partition A .

Figure 2: Spectral clustering algorithm

First, all eigenvalues and their corresponding eigenvectors of the similarity matrix are calculated. The eigenvectors $v_i \in \mathbb{R}^P$ can be sorted according to their eigenvalues in descending order, such that $v_0, v_1, \dots, v_r, \dots, v_P$ with $\lambda(v_0) = \max_{i \in \{0, \dots, P\}} \lambda_i$. Now we can choose the first R eigenvectors to have R different clusters. Our set of eigenvectors is defined as $V = (v_0, v_1, \dots, v_r) \in \mathbb{R}^{P \times R}$. We can now define a set $U = (u_0, u_1, \dots, u_p) = V^T \in \mathbb{R}^{R \times P}$, whereas u_i corresponds to exactly one data point and is the vector containing the i -th elements of all eigenvectors v_1, \dots, v_r . As a last step, we apply a weighted version of the k -means algorithm [17] by first randomly selecting R vectors m_1, \dots, m_r as mean vectors and then successively assigning the vectors u_1, \dots, u_p and updating m_1, \dots, m_r . After conversion, all vectors assigned to m_j present the j -th cluster.

Bach and Jordan applied spectral clustering for source separation of two-speaker mixtures. The method described in the following is based on their paper “Learning spectral clustering, with application to speech separation” [2].

First, the definition of a similarity matrix W corresponding to the mixture signal $X(t)$ will be derived. As a first step the mixture signal can be converted to (F, T) bins using the short term fourier transform as explained in 2.2.1. The (F, T) are then used to create multiple features, which are based on: *Energy, continuity, common fate cues, pitch estimation, timbre,...* In this way, we create multiple features $s_i(f, t)$ for every data point (f, t) whereas the calculation of $s_i(f, t)$ can make use of neighboring data points.

The features are essentially embeddings making the data points more comparable, so that we can define a similarity matrix corresponding to k features $W_k(i, j) = f_k(s_i, s_j)$ choosing appropriate functions f_k . The final similarity matrix consists of a weighed combination of each similarity matrix $W = W_1^{\alpha_1} \odot W_2^{\alpha_2} \odot \dots \odot W_k^{\alpha_k}$ with \odot being the hadamard product. The similarity matrix can then be inserted into our previously explained spectral clustering algorithm 2 and the two clusters can be defined resulting in a neat spectrogram for every speaker. The conversion of the data points to a combined similarity matrix works like front-end processor.

The only thing left to explain is how to define the weight vector $\alpha \in \mathbb{R}^K$. Having N labeled data sets with the (F, T) bins corresponding to one of two speakers so that every data set has a reference partition E'_n , a partition $E_n(\alpha) = \text{K-Means}(W_n(\alpha))$ can be derived. The *k-means* algorithm is applied to the similarity matrix depending on α and the previously defined features as well as the *k-means* algorithm. Finally, a cost function $\mathbb{L} = \sum_i^N H(E'_n, E_n(\alpha))$ is used to find the α minimizing the \mathbb{L} .

Even though spectral clustering can produce very good results on two-speaker separation [2] it has its limitations. Already for seconds of speech sampled at 5.5 kHz leads to 22000 data points and thus, a similarity matrix having 22000×22000 entries not even applying the short term fourier transform. It is obvious that eigenvalue and eigenvector calculations become very expensive. A remedy is applying low rank matrix approximations such as the Nystrom approximation to perform eigenvalue and eigenvector on a matrix that is reduced from $N \times N$ to $N \times D$. This allows for significant performance improvement and is especially useful on sparse matrices [2].

Additionally, the algorithm relies on “handcrafted features” and uses only shallow learning for the similarity matrices by essentially applying linear regression to optimize the scaling vector α . More speakers means more overlapping speakers per (F, T) bin so that features taking into account multiple data points for calculation contain multiple speakers making it pointless to cluster them. A speaker segmentation would be needed before-hand to create useful features, which would miss the point of creating features to perform the segmentation. Deep clustering as explained in 4.1 tries to solve this problem.

4 Deep clustering methods

In this section, two state-of-the-art approaches using artificial neural networks for the source separation problem will be discussed. The first approach, called “Deep clustering” ?? is very similar to *spectral clustering* as explained in the section before and works in the frequency domain of the signal. the second approach, called “TasNet” works in the time domain and will be discussed in detail in 4.2. Finally, the subsection introduces a metric for performance measurement and compares the two methods.

When applying artificial neural networks to the source separation problem there are two common problems that occur.

- The permutation problem
- The dimension output problem

To best understand the permutation problem of artificial neural networks, let us go through an example. An artificial network for two-speaker source separation that takes $n(F, T)$ bins corresponding to the mixed signal $X(t)$ as an input in the first layer would need an output layer of size $2 \times n$ with the first half corresponding to the first speaker and the second half to the second speaker. Having three two-speaker datasets, for example (A,B), (A,C) and (B,C) to train such a network the following problem occurs. First the network will learn to activate the (F, T) corresponding to speaker A on the left half of the output layer and the output for speaker B on the right half. Second, the network learns to assign speaker A again to the left half and speaker C to the right half. Third, when training on the dataset of speaker B and C, the network will try to assign the right half to both speaker B and C [6]. These constant “mis-learnings” will it most probably impossible for the learning process to converge.

The dimension output problem refers to the static input and output dimension of normal feed forward neural networks. Having defined a output dimension corresponding to two speakers for example the network will learn to separate only two-speaker signal mixtures making it impossible to separate signal mixtures containing more than two speakers.

The explanation of the two methods will be structured by answering the following questions.

1. What is the idea of the method?
2. How is the system trained?
3. How is the trained system applied to new data?
4. How does the method tackle the two problems described above?

4.1 Deep clustering method - frequency domain audio separation based methods

Deep clustering was introduced by John R.Hershey and co. [10]. It extends the application of spectral clustering by applying artificial neural networks to model an embedding V replacing the similarity matrix W as described in 3.2. The artificial neural network is represented by $f_\theta(S) = V$ with

- θ being the parameter space of the artificial neural network
- S with $s = s_{t,f} = s_n \in S \in \mathbb{R}^N, N = T \times F$
- V with $v = v_n \in \mathbb{R}^K$ and $v_n \in V \in \mathbb{R}^{N \times K}$ and $|v_n|^2 = 1$

S is the spectrogram or a feature extracted spectrogram of the mixed signal $X(t)$. V is a K -dimensional embedding, with each v_n representing a normalized embedding for s_n (one bin of the (T,F) spectrogram). Clustering of v_n using the k-means algorithm

then results in binary masks that can be applied to the (T,F) spectrogram to retrieve the separated signals for every speaker. It is obvious that much more complex structures and relations can be learned in *deep clustering* than in *spectral clustering*.

To train the network, the following cost function is applied to θ . For practical reasons we will write $V(\theta)$ simply as V , but keep in mind that V depends on θ .

$$C(\theta) = |VV^T - YY^T|_F^2 \quad (1)$$

with $C(\theta)$ the cost function and

- $VV^T \in \mathbb{R}^{N \times N}$ being the similarity matrix and v
- $Y \in \mathbb{R}^{N \times C}$ with C being the number of different speakers. $y_{n,i} = 1$ if speaker i is dominating in the (T,F) bin n and $y_{n,i} = 0$ otherwise. Thus y_n presents a one-hot encoded vector. Thus $YY^T(n_1, n_2) = 1$ only if $y_{n_1,i} = y_{n_2,i}, \forall i \in 1, \dots, c$ and 0 otherwise.
- F being the Frobenius norm with $|A|_F^2 = \sum_{i,j} A(i,j)^2$

Having the optimal source separation Y^* we can derive Y using a one-hot encoding replacing dominante values with 1 and the rest with 0. Equation 1 can be rewritten.

$$\begin{aligned} C(\theta) &= |VV^T - YY^T|_F^2 = \sum_{i,j} (v_i^T v_j - y_i^T y_j)^2 \\ &= \sum_{i,j; y_i=y_j} (v_i^T v_j - 1)^2 + \sum_{i,j; y_i \neq y_j} (v_i^T v_j)^2 \\ &= \sum_{i,j; y_i=y_j} (v_i^T v_j)^2 - 2(v_i^T v_j) + 1 + \sum_{i,j} (v_i^T v_j)^2 - \sum_{i,j; y_i=y_j} (v_i^T v_j)^2 \\ &= \sum_{i,j; y_i=y_j} (-2v_i^T v_j + 1) + \sum_{i,j} (v_i^T v_j)^2 \end{aligned} \quad (2)$$

Using the “parallelogram law”: $u^T v = \frac{1}{2}(|u|^2 + |v|^2 - |u - v|^2)$ with u, v being vectors, we get

$$\begin{aligned} C(\theta) &= \sum_{i,j; y_i=y_j} (-(|v_j|^2 + |v_i|^2 - |v_i - v_j|^2) + 1) + \sum_{i,j} (v_i^T v_j)^2 \\ &= \sum_{i,j; y_i=y_j} (|v_i - v_j|^2 - 1) + \sum_{i,j} (v_i^T v_j)^2 \\ &= \sum_{i,j; y_i=y_j} |v_i - v_j|^2 + \sum_{i,j} (v_i^T v_j)^2 - N \end{aligned} \quad (3)$$

We derived an equation that allows for a more intuitive representation: For $y_i = y_j$ the respective embeddings v_i and v_j should be as close together as possible when trying to minimize the first addend. The second addend and third addend can be seen as a regularizer preventing trivial solutions. As a conclusion, the embedding V is very similar to the low rank approximation matrix used in spectral clustering.

To optimize computation performance the expensive $N \times N$ matrix calculation in equation ?? can be replaced by the less expensive $K \times K$ with K being the embedding dimension of V which is normally magnitudes smaller than N . We arrive at:

$$C(\theta) = |V^T V|_F^2 - 2|V^T Y| + |Y^T Y| \quad (4)$$

Finally, the derivations $\frac{\partial C(\theta)}{\partial \theta}$ can be split into $\frac{\partial C(\theta)}{\partial V} \frac{\partial V}{\partial \theta}$. While the calculation of $\frac{\partial V}{\partial \theta}$ depends heavily on the network architecture and can be quite expensive using, for example long short term memory units, as in [6], it is easy to calculate:

$$\frac{\partial C(\theta)}{\partial V} = 4V(V^T V) - 4Y(Y^T V)$$

The parentheses in the above equation make sure that the computational less expensive matrix operations are executed.

Having succesfully trained the parameters θ of the neural network, we can now compute V of unseen data X using the short term fourier transform and applying the neural network $V = f_\theta(\text{ShortTermFourierTransform}(X))$. Analog to *spectral clustering*, we use the *k-means* algorithm by minimizing its inference cost with the following equation:

$$\hat{Y} = \operatorname{argmin}_Y V - Y(Y^T Y)^{-1} Y^T V \quad (5)$$

\hat{Y} are used as binary masks to separate the sources. So we can lastly apply the masks $\hat{Y}(c)$ dot-wise to the spectrogram of the mixture signal:

$$S_c = \hat{Y}(c) \odot X$$

Since $(Y^T Y)^{-1} Y^T \in \mathbb{R}^{C \times D}$ presents the means of all clusters, minimizing the cost function 5 is essentially the same as performing k-means clustering on $v_n \in V$ with successive creating of the binary masks from the assignments [6]. In the case that $VV^T = Y^* Y^T$ both cost functions 5, 1 are coincide in their global minimum.

Since training essentially relies on YY^T both the permutation and the dimension output problem are solved. $YY^T(n_1, n_2) = YY^T n_2, n_1 = 1$ if y_{n_1} corresponds to the same class as y_{n_2} and 0 otherwise. So the order of the classes is clearly disregarded since only "same" or "different" classes matters. Also YY^T always has the same dimension $N \times N$ no matter how many speaker are mixed in the signal.

As we will later see in section 4.2.1, results achieved by *clustering* are impressive, there is room of improvement considering that the (F, T) are hard assigned to only one class.

4.2 TasNet - time domain audio separation based methods

Introduce TasNet which does not create masks for each source, but instead use encoder-decoder framework to model directly the signal in the time-domain. Use [12].

4.2.1 Results

Talk about results of TasNet

4.3 Comparison of the two methods

5 Conclusion

Write conclusion using results.

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