Friedrich-Alexander-Universität Erlangen-Nürnberg

FRIEDRICH-ALEXANDER
UNIVERSITÄT
ERLANGEN-NÜRNBERG

Master Thesis

Chord Transition Features for Style Classification of Music Recordings

submitted by

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submitted

October 12, 2018

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Erlangen, 12. Oktober 2018

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Abstract

Chords are a central component of Western music. From a musicological viewpoint, chords and chord progressions carry important information about a piece of music. Estimating chord labels from audio recordings is a well-known problem in the field of Music Information Retrieval. The task of chord recognition consists in computing a sequence of chord labels from a recording. A standard approach uses Hidden Markov Models and the Viterbi algorithm. From the resulting chord label sequence, we can obtain local information about the transitions between subsequent chords. However, using transitions derived from a single "optimal" chord sequence is sometimes problematic. Chord labels are often ambiguous and by deciding on one sequence, we cannot capture this ambiguity. This also affects the chord transitions derived from these labels. In this thesis, we propose a novel type of mid-level features which capture local chord transition probabilities. To this end, we exploit the Baum–Welch algorithm, which typically serves to estimate parameters for the Hidden Markov Model. We then propose a non-linear rescaling technique that we need to display chord transition probabilities. In several experiments, we show the benefits of these mid-level features. We test the algorithm in a chord recognition scenario and show the benefits of soft chord transition features by demonstrating how we can use the features to gain additional musical information about a recording. We then test the mid-level features in a genre classification scenario. Concretely spoken, we deal with a sub-style classification scenario, where we distinguish between four musical periods within Western classical music: Baroque, Classical, Romantic, and Modern. For this problem we use a standard machine learning pipeline. To use our features in a classifier, we apply an aggregation algorithm to obtain piece-level features. We find that our features yield comparable results to state-of-the-art tonal audio features. When we compare the proposed features with the corresponding hard-decision chord transition features, we obtain a consistently better performance.

Zusammenfassung

Akkorde bilden eine wesentliche Komponente westlicher Musik. Aus musikwissenschaftlicher Sicht stellen die verwendeten Akkorde und Akkordfolgen relevante Information über ein Musikstück dar. Im Bereich der Music Information Retrieval sind Verfahren bekannt, mit deren Hilfe Akkorde ausgehend von Musikaufnahmen geschätzt werden können. Ein Akkorderkennungsalgorithmus berechnet eine Akkordfolge für eine Aufnahme, oft mithilfe eines Hidden Markov Models und dem Viterbi-Algorithmus. Auf Basis dieser Folge können wir nun Aussagen über die Übergänge aufeinanderfolgender Akkorde treffen. Es ist jedoch oft problematisch, eine eindeutige, „optimale“Akkordfolge zur Beschreibung von Akkordübergängen zu bestimmen. Akkorde können oft nicht zweifelsfrei zugeordnet werden und durch die Entscheidung auf eine Sequenz ist es nicht möglich, diese Mehrdeutigkeit zu erfassen. Davon werden auch abgeleitete Messungen zu Akkordübergängen beeinflusst. In dieser Arbeit stellen wir eine neue Art von „weichen“Akkordübergangsmerkmalen vor, die auf Schätzungen lokaler Übergangswahrscheinlichkeiten beruhen. Hierzu nutzen wir den Baum-Welch-Algorithmus, der typischerweise zur Bestimmung von Modellparametern von Hidden Markov Models herangezogen wird. Zudem stellen wir eine neue nichtlineare Umskalierungsmethode vor, die wir benötigen um Akkordübergangswahrscheinlichkeiten darzustellen. Wir zeigen in den nachfolgenden Experimenten die Vorteile, die sich aus solchen „weichen“Merkmalen ergeben. Wir testen unseren Algorithmus auch in einem Akkorderkennungsszenario und zeigen, wie wir mithilfe von „weichen“Merkmalen zusätzliche Informationen über ein Musikstück gewinnen können. Als ein Hauptbeitrag dieser Arbeit zeigen wir die Qualität der Merkmale indem wir sie auf ein Klassifizierungsproblem anwenden. In unserer speziellen Aufgabenstellung klassifizieren wir Aufnahmen klassischer Musik in vier Epochen: Barock, Klassik, Romantik und Moderne. Wir verwenden hierzu Standardalgorithmen für maschinelles Lernen. Um unsere Merkmale nutzen zu können, verwenden wir einen Aggregationsalgorithmus, mit dem wir die lokalen Merkmale zu globalen Merkmalen, die eine ganze Aufnahme beschreiben zusammenfassen. Unsere Merkmale liefern vergleichbare Ergebnisse zu anderen tonalen Merkmalen aus der Literatur. Unsere Experimente zeigen, dass die neuen, „weichen“Merkmale durchgehend bessere Ergebnisse erzielen als die konventionellen, „harten“Merkmale.

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Chapter 1

Introduction

The notion of chords is central when analyzing Western music. A chord describes several notes which are played simultaneously, we can see it as a vertical component of music [21]. Knowledge of the chord sequence of a piece of music is useful in a number of scenarios: There are applications in music processing tasks such as music structure analysis or music retrieval [24]. Of particular interest in this thesis are chord progressions, which carry important information about a piece of music [16]. We can gain knowledge about chord transitions by analyzing symbolic data, such as MIDI files or annotated chord labels [27]. If there is no annotation available, we need to use a chord recognition approach. In this thesis we focus on the case that there is only an audio recording available. In this context, chord recognition describes the estimation of a chord sequence from a given recording. This is a well-known task in the field of Music Information Retrieval (MIR) [9, 24]. There are various approaches available to estimate a chord sequence [9], the most common of which is based on Hidden Markov Models. Sheh and Ellis [36] proposed to use the Viterbi algorithm [43] to estimate a chord sequence. The resulting unique chord sequence is optimal in a certain way. We can use this sequence to obtain local information about transitions between subsequent chords [45].

However, using a chord sequence to derive chord transitions is sometimes problematic. Even though the chord sequence is "optimal", it still might not always be correct from a musicological perspective. Since every transition depends on two chords, an error in just one of the chords is sufficient to yield an error in the estimated transition. Since every chord label influences two chord transitions, one chord-estimation error results in two chord transition errors. Also, chord labels are often ambiguous [24, Section 5.2.2]. Deciding on one chord sequence ignores this ambiguity, which also affects the estimated chord transitions. In this thesis we propose a novel approach of estimating chord transitions in a soft way to solve this problem. We propose a method which is based on the Baum–Welch algorithm [3] instead of the commonly used Viterbi algorithm. We use results within this algorithm, based on the Forward-Backward algorithm [3] to obtain probabilities

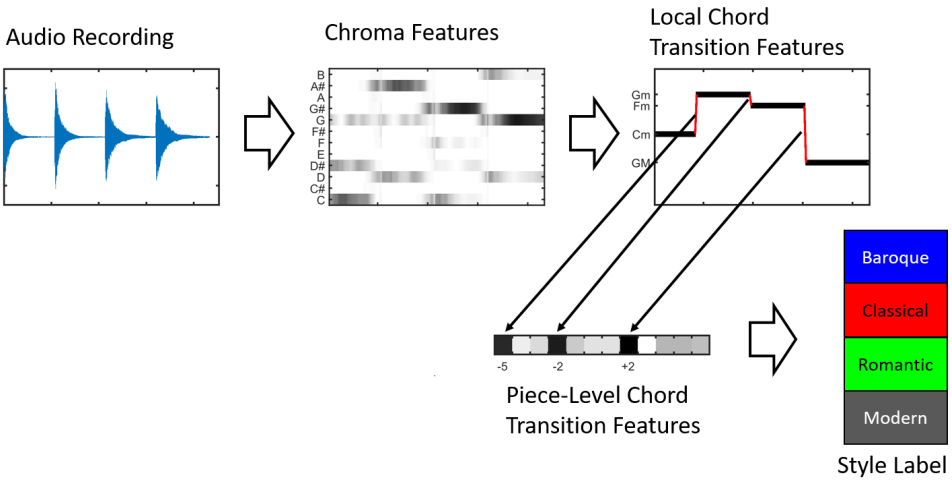


Figure 1.1: Processing pipeline from the audio recording to the estimated style label.

for all chord transitions for each frame. As main contribution we propose such mid-level chord transition features and demonstrate their performance in several experiments. First, we show that conventional visualization techniques are not sufficient to properly show the chord transition probabilities. We therefore propose a novel, non-linear rescaling technique to visually enhance the plotted transition probabilities. We demonstrate how soft chord transition estimation can be useful to gain further musical knowledge about a recording, using a choral by J.S. Bach as an example. We then use our features to perform sub-style classification. In our particular task—which is related to the widely-known problem of genre classification [30, 39, 41]—we classify different recordings of Western classical music from four different musical periods: Baroque, Classical, Romantic, and Modern. To this end, we employ an aggregation algorithm to transform the local chord transition features into piece-level features. We then use these features in a standard machine learning pipeline. While the machine learning techniques are also important in this thesis, the main contributions are the proposed features and the systematic evaluation in a style classification scenario.

Within the larger context of piece-level audio features, we can categorize our proposed features as mid-level tonal features. Tonal features are musically motivated and are often based on pitch class energy distributions [17]. We can further distinguish between high-level features and mid-level features [45]. High-level features directly describe concepts which are interpretable by humans, for example the number of a certain type of chord transitions. Therefore we count chord transition features which are based on a single unique chord sequence among the high-level features. Mid-level features, on the other hand, describe those concepts in a less obvious way, for example by describing soft transition probabilities. Therefore, our proposed features belong to the mid-level features.

One main contribution of this thesis is the application in a style-classification scenario. Figure 1.1 gives an overview of the pipeline we use. At first, we transform the waveform into the musically more informative chroma features. We then employ either the Baum–Welch algorithm or the Viterbi algorithm to derive local chord transition features. Afterwards, we aggregate the local features into one piece-level feature vector. We use this feature vector in a classifier to estimate the musical style. This thesis explains every component of this pipeline, but we focus on the estimation of local chord transition features.

The remainder of the thesis is structured as follows:

In Chapter 2, we introduce basic musical concepts, both on a musical and on a technical level. We define basic notions on which the thesis is built. Afterwards, we introduce the task of chord recognition and discuss basic algorithms to solve this task. To that end, we introduce Hidden Markov Models—a concept we use to model chords as hidden internal states of a recording.

Chapter 3 covers the generation of chord transition features. This is one central contribution of this thesis. We describe the feature generation using the Viterbi algorithm for computing hard-decision features and the Baum–Welch algorithm for computing soft-decision features. We introduce a novel visualization technique for displaying chord transition probabilities employing a special rescaling procedure. Afterwards, we first evaluate both algorithms on a chord recognition task. We examine the effect of the self-transition probability—a central parameter in our model—by performing systematic experiments on 180 Beatles songs. We discuss the resulting local chord transition features, before we introduce and discuss a feature aggregation algorithm to obtain piece-level features.

In Chapter 4, we introduce the problem of style classification and give an overview over the dataset we use for our experiments. We then give a brief introduction to some machine learning concepts we need for this thesis. Finally, we give an overview about related work on this topic.

In Chapter 5, we present experiments for style classification and discuss the results. We compare our features against state-of-the-art tonal audio features evaluated in related work. We focus on the comparison between hard-decision and soft-decision chord transition features in different scenarios. We further investigate possibilities of combining our features with state-of-the-art tonal features. We test chord transitions features in different settings regarding the training and test set to demonstrate the so-called “album effect”. As a final experiment, we validate our results by performing a cross-version experiment.

The thesis ends with a summary of the obtained results and a brief discussion of open issues which could be examined in future work.

Chapter 2

Background: Chord Recognition

This chapter gives a brief introduction to the broad field of musical chord recognition. We start with some musical foundations on a symbolic level. Afterwards, we introduce chroma features, a common feature representation for the task of chord recognition [14]. Building on this, we discuss the straightforward approach of template-based chord recognition, before moving on to the more sophisticated Hidden Markov Model (HMM) based approach. We use a recording of the choral "Durch dein Gefängnis, Gottes Sohn" by Johann Sebastian Bach (BWV 245, No. 22) performed by the Choir of Kings College Cambridge as an example throughout this chapter.

2.1 Symbolic Music Representations

In this section, we discuss musical concepts using symbolic representations. In the first part, we cover the basics concerning notes and pitches before moving on to intervals and chords.

2.1.1 Notes and Pitches

The following is a summary of [24, Section 1.1]. The most basic element of Western music notation is a *note*. A note usually corresponds to a tone of a certain fundamental frequency or *pitch* played for a certain duration. Pitch is a perceptional property, which effectively describes how "high" a tone is. Since the human perception of frequency is logarithmic, it is common practice to give pitch on a logarithmic scale. A musical composition consists of many notes ordered in a specific way. We can loosely define *melody* as a temporal progression of notes and *harmony* as several notes being played at the same time, a kind of vertical component of music [21]. It has been found that notes whose frequencies have the ratio of an integer power of two are perceived as similar. Therefore, two notes with this property are summarized to one *pitch*

(a) (b)

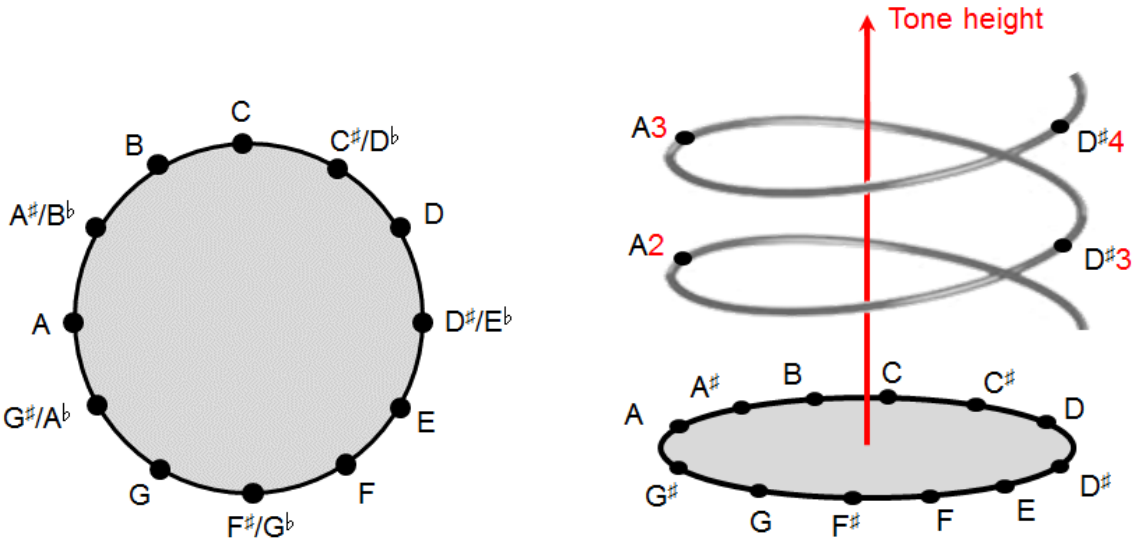


Figure 2.1: (a) Chromatic circle and (b) Shepard's helix of pitch perception. Both images are taken from [24, Figure 1.3]

class. With this we can define an *octave* as the distance between two notes with a pitch ratio of two. In general, the pitch scale is continuous but in Western music, it is common to discretize the pitch on a logarithmic frequency axis such that each octave splits into twelve intervals of the same size. The result is called the twelve-tone equal-tempered scale. The distance between two adjacent notes is called a *semitone*. Two notes that are one semitone apart have a frequency ratio of $\sqrt[12]{2} \approx 1.059$. Within the equal tempered scale, each tone can be assigned a pitch class and an octave. In the international scientific pitch notation, each pitch class is assigned a letter from A to G and possibly, an accidental (♯ or ♭) shifting the tone one up or down a semitone. At this point, we assume that the reader is familiar with the basic concept of music notation. An important observation here is the cyclic nature of pitch classes [37]. By repeatedly increasing the pitch class by one step, we will arrive at the original pitch class again after 12 steps. This is nicely visualized by the so-called chromatic circle (Figure 2.1a). By adding the octave information, we obtain the so-called Shepard's helix of pitch perception (Figure 2.1b), in which each note is ordered on an ascending helix, where the angle represents the pitch class and the elevation the height of the tone. These representations show another important concept, *enharmonic equivalence*. Within this concept, it is not possible to distinguish pairs like $D^{\flat}$ and $C^{\sharp}$. Notes with different letter names but representing the same pitch-class are called enharmonically equivalent.

Distance (Semitones)	Name	Example	Frequency Ratio	Approximate Ratio
0	(Perfect) Unison	C-C	1.000	1 : 1
1	Minor second	C-D$^\flat$	1.059	16 : 15
2	Major second	C-D	1.122	9 : 8
3	Minor third	C-E$^\flat$	1.189	6 : 5
4	Major third	C-E	1.260	5 : 4
5	(Perfect) fourth	C-F	1.334	4 : 3
6	Tritone	C-F$^\sharp$	1.414	45 : 32
7	(Perfect) fifth	C-G	1.498	3 : 2
8	Minor sixth	C-A$^\flat$	1.587	8 : 5
9	Major sixth	C-A	1.681	5 : 3
10	Minor seventh	C-B$^\flat$	1.781	9 : 5
11	Major seventh	C-B	1.888	15 : 8
12	(Perfect) Octave	C-C	2.000	2 : 1

Table 2.1: Names of the intervals ordered by their distance in semitones. We give examples for each interval with root note C. Also we show the frequency ratio in the equal-tempered scale and an approximate frequency ratio [24, Table 5.3].

2.1.2 Chords and Intervals

Following [24, Section 5.1] and [45, Section 2.6.1], this section presents information and definitions for the concepts of chords and intervals. A *chord* is loosely defined as a set of notes played at the same time.

Before presenting different kinds of chords, we first introduce the concept of an *interval*. We already mentioned an important interval when we defined the octave as distance between two notes with a pitch ratio of two. In fact, each interval is the distance between two notes with a certain pitch ratio. In this context, "distance" refers to the difference of the fundamental frequencies on a logarithmic scale. On the equal-tempered scale, those ratios can assume irrational values. However, they can be approximated by fractions. Intervals with fractions of small numbers sound rather pleasant. Intervals described by more complicated fractions sound unpleasant. This is called consonance and dissonance, respectively. Table 2.1 lists all intervals of the chromatic scale up to an octave with the corresponding pitch ratios and an approximation as a fraction. From this table, we see that the fourth and fifth intervals are particularly consonant. Also, the thirds and sixths are perceived as consonant, but the higher integers in the fraction indicate a lower level of consonance. The other chords are considered to be dissonant [21].

With this in mind, we can be loosely define a chord as several notes being played simultaneously. The first step towards classifying chords is to divide them by the number of notes. In this thesis, we only consider *triads* (three notes) and *tetrads* (four notes). Each note has a so-called *root*

CM Cm C+ C°

Figure 2.2: Examples of four triad types in root position.

note, which is the basis of this chord. Often but not always, the root note has the lowest pitch of the chord. In this case we speak of a chord in *root position*.

The most important triads in Western music are tertian chords, which consist of notes that can be stacked in thirds [21], meaning that the interval between two adjacent notes is either a major or a minor third. The most common triads are the *major* and *minor* triads. The major triad consists of a major third above the root note and a minor third on top of it. The minor triad consists of a minor third and a major third on top. In both cases, the interval between the lowest and highest note is a perfect fifth, giving the chords a stable quality. We denote a major triad with its root pitch class followed by a capital M. For example, the C major triad is written as CM. The we denote minor triads with a lower-case m, so we abbreviate the C minor triad by Cm. Figure 2.2 gives an example for both triad types.

A way of mathematically describing a chord is to list the pitch class of all involved notes in a pitch class set S. We number the pitch classes from 0 to 11 starting with C and ascending in semitone steps. We can write the pitch class set of CM, which consists of C, E and G, as:

$$S_{\mathrm{CM}} = \{0, 4, 7\} \tag{2.1}$$

Another useful representation are chord template vectors. In these twelve-dimensional vectors, each entry corresponds to one of the pitch classes. If the pitch class is active in the chord, the corresponding entry is set to 1 otherwise the entry is 0. As an example, we can write the template vectors of CM and Cm as:

$$\begin{aligned} \mathbf{t}^{\mathrm{CM}} &= (1,0,0,0,1,0,0,1,0,0,0,0)^{\intercal} \\ \mathbf{t}^{\mathrm{Cm}} &= (1,0,0,1,0,0,0,1,0,0,0,0)^{\intercal} \end{aligned} \tag{2.2}$$

There are two other possibilities to stack two third intervals by using two major thirds or two minor thirds. Those chords are called the *augmented* and *diminished* triad, respectively. Note that the interval between the highest and lowest is no longer a perfect interval. The augmented chord has a minor sixth, also called augmented fifth interval. The diminished chord has a tritone or diminished fifth interval, which is particularly dissonant. The chords are denoted by "+" for augmented chords and "°" for diminished chords. We show note examples of augmented

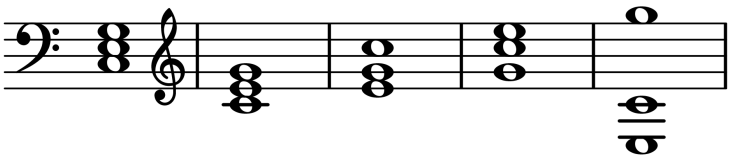


Figure 2.3: The CM triad in several variations. All shown chords comprise the pitch classes C, E and G, but in different octaves. Even though the musical impressions may be quite different they all correspond to the same template vector.

and diminished triads in Figure 2.2 and give the template vectors with the root note C in the following:

$$\begin{aligned} \mathbf{t}^{\mathrm{C}+} &= (1,0,0,0,1,0,0,0,1,0,0,0)^\intercal \\ \mathbf{t}^{\mathrm{C}^\circ} &= (1,0,0,1,0,0,1,0,0,0,0,0)^\intercal \end{aligned} \tag{2.3}$$

The template vectors for different root notes can be obtained by cyclically shifting the template vector of C. The number of shifts equals the distance of the respective pitch classes. For example, the template vector for Gm can be written as:

$$\mathbf{t}^{\mathrm{Gm}} = (0,0,1,0,0,0,0,\mathbf{1},0,0,1,0)^\intercal\,. \tag{2.4}$$

For clarity, the entry representing the root note is printed in bold face. By describing a chord only as a set of pitch classes, there are several different chords corresponding to the same representation. First, since the octave information is lost, chords played in different octaves are mapped to the same template vector. Furthermore, inversions—in which the root note is not the one with the lowest pitch—become indistinguishable. The notes of a chord may even be distributed over several octaves. Figure 2.3 shows several chords which are all mapped to the same template vector.

This property is particularly interesting when it comes to symmetric chords, for example the augmented chord. Without respecting the octave information, it is impossible to distinguish more than four root notes. Since C+, E+ and G♯+ all comprise the pitch classes C, E and G♯ they have the same representation, because of the aforementioned enharmonic equivalence. This will become an issue later on.

This concept can be extended to tetrads or—more specifically—seventh chords, which consist of a triad and a seventh interval. The chords we use in this thesis consist of three stacked major or minor thirds. There are eight possibilities, but stacking three major thirds leads to a repetition of the root pitch class when respecting enharmonic equivalence. Therefore, the chord is an augmented chord. Hence, seven types of seventh chords remain.[1] In Table 2.2, we present the

[1] There are actually more seventh chords than we described here. However, this is the set we are working with.

Name	Notation	Template Vector
Major seventh	$\mathrm{CM}^{\mathrm{maj7}}$	$\mathbf{t}^{\mathrm{CM}^{\mathrm{maj7}}} = [1,0,0,0,1,0,0,1,0,0,0,1]^\intercal$
Minor seventh	Cm^{7}	$\mathbf{t}^{\mathrm{Cm}^{7}} = [1,0,0,1,0,0,0,1,0,0,1,0]^\intercal$
Dominant seventh	CM^{7}	$\mathbf{t}^{\mathrm{CM}^{7}} = [1,0,0,0,1,0,0,1,0,0,1,0]^\intercal$
Diminished seventh	$\mathrm{C}^{\circ 7}$	$\mathbf{t}^{\mathrm{C}^{\circ 7}} = [1,0,0,1,0,0,1,0,0,1,0,0]^\intercal$
Half-dim. seventh	$\mathrm{C}^{\emptyset 7}$	$\mathbf{t}^{\mathrm{C}^{\emptyset 7}} = [1,0,0,1,0,0,1,0,0,0,1,0]^\intercal$
Min. maj. seventh	$\mathrm{Cm}^{\mathrm{maj7}}$	$\mathbf{t}^{\mathrm{Cm}^{\mathrm{maj7}}} = [1,0,0,1,0,0,0,1,0,0,0,1]^\intercal$
Aug. maj. seventh	$\mathrm{C}+^{7}$	$\mathbf{t}^{\mathrm{C}+^{7}} = [1,0,0,0,1,0,0,0,1,0,0,1]^\intercal$

Table 2.2: Template vectors of the different seventh chords we use in this thesis.

template vectors of all used seventh chords for reference. When inspecting the diminished seventh chord, we can see that—similarly to the augmented triad—the template vector is symmetric. In this case, there are only three distinguishable root notes.

In this section, we gave an overview over the basic building blocks of Western music. We limited ourselves to concepts which are necessary to follow the technical content of the thesis. For a more detailed discussion of musical theory, we refer to [45, Chapter 2].

2.2 Feature Representation

In the previous section, we discussed some musical concepts on a symbolic level. Now, we introduce several ways to represent music recordings in a technical context. This part is based on [24, Chapter 3.1] with a brief part taken from [24, Chapter 1.3.1].

An audio recording captures the relative air pressure as measured by a microphone which can be reproduced by a loudspeaker. Typically, 44100 or 22050 samples are recorded per second. A sampling rate F_s of more than 40000 Hz is sufficient to represent all frequencies audible for humans. We now want to transform the waveform into a musically meaningful representation with a lower sampling rate. As a first step, most MIR approaches apply the short-time Fourier transform (STFT) in order to transform the waveform into a so called *spectrogram*, which is used to analyze the predominant frequencies. An extensive discussion of the different kinds of Fourier transforms and the STFT can be found in [24, Chapter 2]. The STFT splits the signal into small overlapping windows and transforms each window into the frequency domain. The resulting amplitude vectors are then concatenated to obtain a two dimensional representation $\mathcal{X}(n,k) \in \mathbb{R}^{N\times K}$, with N and K as the number of time frames and frequency bins, respectively. This contains information about the energy of the signal at time frame $n \in [1:N] := \{1,2,\ldots,N\}$, and frequency index $k \in [0:K]$. Each value of k corresponds to a frequency $F_\mathrm{coeff}(k) = \frac{kF_\mathrm{s}}{N_w}$ with the STFT window size N_w.

As described in the previous section, the human perception of frequency is logarithmic. Therefore, we order the frequencies logarithmically. To get a representation close to the musical context, we bin the frequencies such that they match the pitch scale discussed in Section 2.1. We obtain the so-called log-frequency spectrogram $\mathcal{Y}_{\mathrm{LF}} \in \mathbb{R}^{N \times 128}$ by summing up all frequency bins close to the mid frequency of each pitch $p \in [0:127]$

$$\mathcal{Y}_{\mathrm{LF}}(n,p) = \sum_{k \in P_{\mathrm{p}}(p)} |\mathcal{X}(n,k)|^2, \tag{2.5}$$

where $P(p)$ is the set of all frequency indices k belonging to the pitch p

$$P_{\mathrm{p}}(p) = \{k : F_{\mathrm{pitch}}(p-0.5) \leq F_{\mathrm{coeff}}(k) < F_{\mathrm{pitch}}(p+0.5)\}, \tag{2.6}$$

with the center frequency of pitch p

$$F_{\mathrm{pitch}}(p) = 2^{(p-69)/12} \cdot 440\,\mathrm{Hz}. \tag{2.7}$$

With this definition the note A4 (concert pitch) is assigned to pitch number 69 and the note C4 (middle C) corresponds to pitch number 60. The resulting spectrogram $\mathcal{Y}_{\mathrm{LF}}(n,p)$ captures the energy of the pitch with index p at time index n.

In many cases—as in chord recognition—we are not interested in the octave of a note. Instead the pitch class is carrying the relevant information. Therefore, we convert the log-frequency spectrogram into a pitch class representation, also known as a *chromagram* [44]. The idea behind the chromagram $\mathcal{C} \in \mathbb{R}^{N \times 12}$ is straightforward as we sum up the energies of all pitches belonging to the same pitch class. Those are all the pitches with indexes that are congruent modulo 12 to each other. We define:

$$\mathcal{C}(n,c) = \sum_{p \in P_{\mathrm{c}}(c)} \mathcal{Y}_{\mathrm{LF}}(n,p), \tag{2.8}$$

with $P_{\mathrm{c}}(c) = \{p \in [0:127] \mid p \bmod 12 = c\}$ and $c \in [0:11]$. The matrix $\mathcal{C}$ can be interpreted as a temporal sequence of 12-dimensional vectors $\boldsymbol{x}$. Those vectors are called *chroma vectors* or *chroma features*. In addition, the chroma features are usually normalized either with respect to the ℓ_1 or the ℓ_2 norm. In this thesis, we use the ℓ_1 norm. This way the features are invariant to dynamic changes. To avoid division by zero in frames of silence, which leads to degenerated features, we use a thresholding procedure. If the total energy of a frame is below a certain threshold, all entries of the vector are set to $\frac{1}{12}$. The feature space $\mathcal{F}$ of the chroma features therefore comprises all twelve-dimensional vectors with positive entries that sum up to 1:

$$\mathcal{F} = \left\{\boldsymbol{x} \in \mathbb{R}^{12}_{\geq 0} : \|\boldsymbol{x}\|_1 = 1\right\}, \tag{2.9}$$

with the set of all non-negative real numbers $\mathbb{R}_{\geq 0}$.

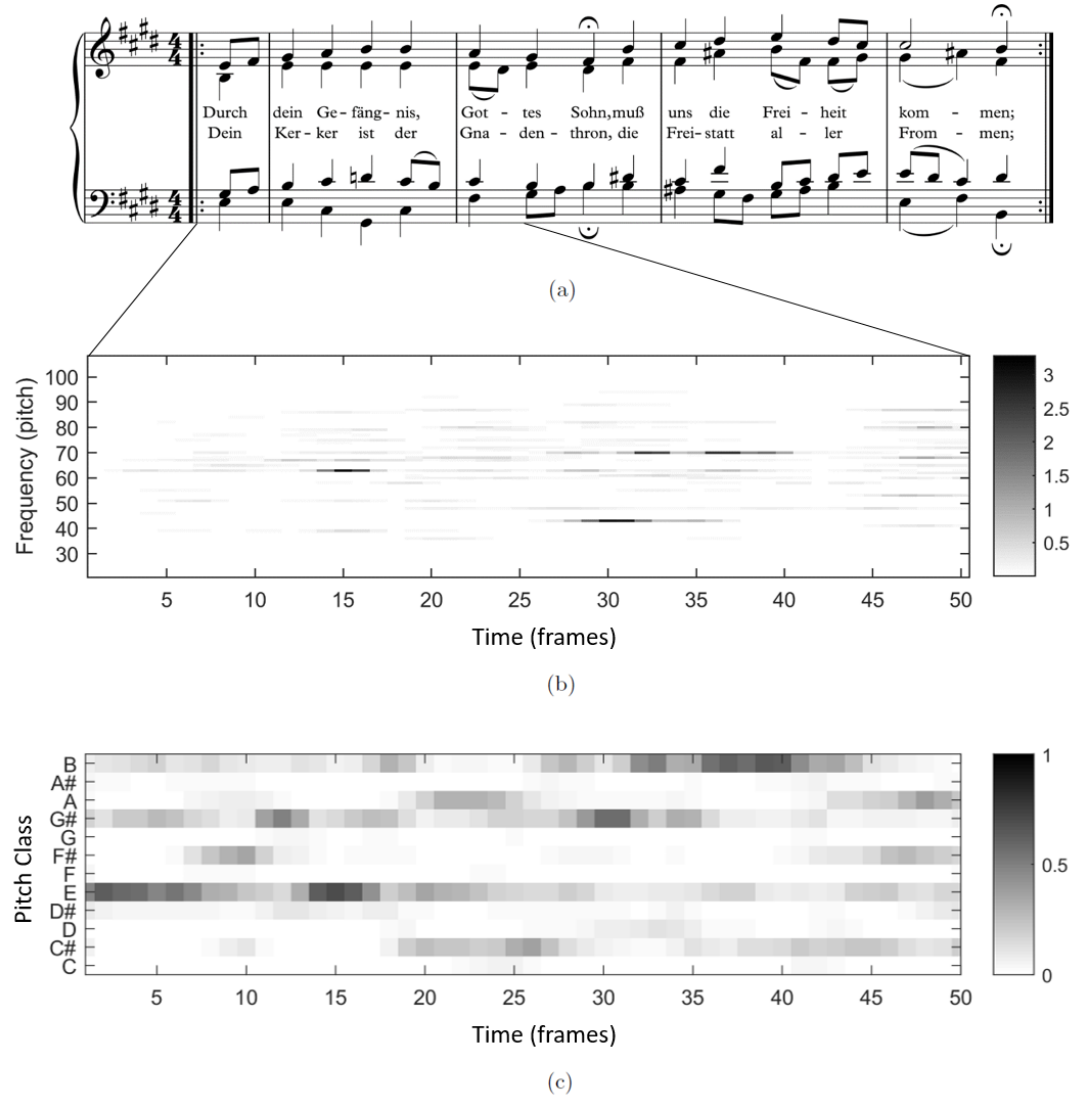


Figure 2.4: Log-frequency spectrogram (a) and chromagram (b) for the first 50 frames of the Bach choral example

In Figure 2.4, we show the log-frequency spectrogram and the chroma features of the Bach choral example. We see that a high value in the spectrogram leads to a high value in the chromagram. However, we can also see high values in the chromagram, where there is apparent silence in the spectrogram (e.g. in the beginning of the recording). This is due to the normalization. This example demonstrates the invariance with respect to changing dynamics. We compute those features using the publicly available Chroma Toolbox[2] [25]. The method for obtaining the log-frequency spectrogram differs from the introduced STFT method by using elliptic filters to directly compute the pitch energies.

[2] `www.audiolabs-erlangen.de/resources/MIR/chromatoolbox`

There are several possibilities how to further enhance the chroma features. One problem of the described chroma features is that different instruments generate a high dynamic range in log-frequency spectrograms. If we generate chroma features from such a spectrogram, we obtain very different chroma feature for different recordings. Especially if we use the features in classifiers this difference will yield problems [9]. A common method to overcome this problem is the logarithmic compression, which effectively limits the dynamic range of the spectrogram. There are several approaches for logarithmic compression found in [9, 24, 26]. Cho and Bello [9] define logarithmic compression as follows:

$$\mathcal{Y}_{\mathrm{LF}}^{(\eta)}(n,p) = \log\left(1 + \eta \cdot \frac{\mathcal{Y}_{\mathrm{LF}}(n,p)}{\max_{\tilde{p}} \mathcal{Y}_{\mathrm{LF}}(n,\tilde{p})}\right), \tag{2.10}$$

with the parameter η, which is typically chosen between 100 and 10000. The obtained spectrogram is then folded into chroma vectors as described in Eq. 2.8.

Another problem concerns with *overtones*. A note played by a real world instrument produces the fundamental frequency of the tone and additional overtones. The frequencies of the overtones are integer multiples of the fundamental frequency. For example a concert A played on a piano comprises the frequencies 440 Hz, 880 Hz, 1320 Hz, 1760 Hz, 2200 Hz, ... These frequencies approximately correspond to the pitch classes A, A, E, A and C$^\sharp$, respectively[3]. The spectral power of each overtone strongly depends on the instrument. In fact, this is one of the characteristics which give each instrument its unique sound. However, this poses an issue if we are only interested in the fundamental frequencies. From the example mentioned above, we see that a played note contributes to more than one pitch class, which makes it considerably harder to get information about the musical content.

An easy approach is octave suppression. Here, a logarithmic frequency spectrogram is multiplied with a window function, reducing the energies in the high and low pitches. The underlying assumption is that the fundamental frequencies are located in the middle of the spectrum. The high frequencies consist mostly of overtones, whereas the low frequencies usually have a bad resolution, due to properties of the STFT and the logarithmic binning. Also, the low-frequency bins a often distorted by a base drum in pop or rock music [9]. In [9] Cho and Bello use a Gaussian window centered around pitch C4 over the logarithmic frequency axis and compute the octave suppressed log-frequency spectrogram as

$$\mathcal{Y}_{\mathrm{LF}}^{OS}(n,p) = \mathcal{Y}_{\mathrm{LF}}(n,p) \cdot \exp\left(-\frac{(p-60)^2}{2 \cdot 15^2}\right). \tag{2.11}$$

[3]Note that the first few overtones are all part of the AM triad. This explains the pleasant sound of major triads, since all notes are consonant with the natural overtones of the root note.

In [23], Mauch and Dixon propose a more sophisticated approach. The proposed chroma features, called non-negative least squares (NNLS) chroma, are based on a fundamental frequency spectrogram, which contains only the fundamental frequencies without the overtones. This spectrogram is estimated by using an overtone model with geometrically declining amplitudes. The amplitude of the h^{th} overtone relative to the amplitude of the fundamental frequency is modeled as s^h with the spectral shape parameter $0 \leq s \leq 1$. We are then dealing with an optimization problem trying to find the fundamental pitches that—in combination with the overtone model—best explain the measured spectrogram. The resulting spectrogram is transformed into chroma features as before. The concept of NNLS chroma was implemented in a Vamp plug-in[4], which can be used in open source software like Sonic Visualizer[5] or Sonic Annotator[6]. This plug-in also contains a spectral whitening technique, which has a similar effect as logarithmic compression. In [9] Cho and Bello extensively study the different extensions of the chroma features and compare them in the context of chord recognition. In the following, we use NNLS chroma with $s = 0.7$.

2.3 Template-Based Chord Recognition

In the previous sections, we discussed the theoretical foundations for chord recognition. This section describes and discusses a straightforward approach for chord recognition based on *template matching*. This section is based on [24, Chapter 5.2] and [9]. Chroma-based chord recognition was initially introduced by Fujishima [14].

In the previous section, we introduced the chroma feature as a music representation capturing the different pitch classes. We showed that an audio recording can be transformed to a sequence of twelve-dimensional chroma vectors, which capture the energy of each pitch class. Comparing this representation to the chord templates defined in Section 2.1, we find conceptual similarities between both representations. Both vectors describe an energy distribution on the different pitch classes. The similarity of a chroma vector and chord template can be a measure of the probability of the chord in this time frame. We define the template-based similarity measure $s_{\text{t}}\left(\boldsymbol{x}, \mathbf{t}^{\lambda}\right)$ between the chroma vector $\boldsymbol{x}$ and the chord template $\mathbf{t}^{\lambda}$ as

$$s_{\text{t}}\left(\boldsymbol{x}, \mathbf{t}^{\lambda}\right) = \frac{\langle \boldsymbol{x} | \mathbf{t}^{\lambda} \rangle}{\|\boldsymbol{x}\|_1 \cdot \|\mathbf{t}^{\lambda}\|_1}, \tag{2.12}$$

where $\mathbf{t}^{\lambda}$ is the template corresponding to the chord λ and $\langle \cdot | \cdot \rangle$ denotes the inner product of two vectors. $\|\cdot\|_1$ denotes the ℓ_1 norm.

[4] `http://isophonics.net/nnls-chroma`
[5] `http://www.sonicvisualiser.org`
[6] `http://www.vamp-plugins.org/sonic-annotator`

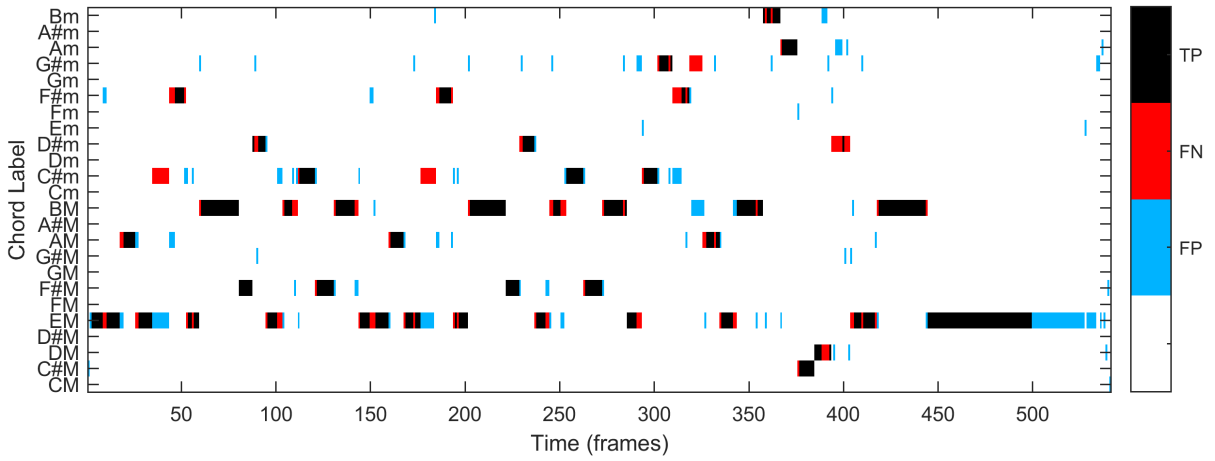


Figure 2.5: Results of template-based chord recognition of the Bach example. The result is color-coded: black represents true positives (TP), red false negatives (FN) and blue false positives (FP). Every color except black and white represents an error.

For estimating a chord label, we pick the chord which maximizes this similarity measure:

$$\hat{\lambda} = \underset{\lambda \in \Lambda}{\operatorname{argmax}}\, s_{\mathrm{t}} \left(\boldsymbol{x}, \mathbf{t}^{\lambda} \right). \tag{2.13}$$

Here Λ denotes a set of chords called a *chord vocabulary*. Usually Λ comprises all 24 major and minor chords [5, 8, 9, 20]. Figure 2.5 shows an example of template-based chord recognition on the Bach example. The plot shows a comparison to the ground truth chords. We obtained the ground truth by manually annotating the recording. For this annotation all chords were mapped on major and minor triads. This is not always justified, especially when dealing with augmented and diminished chords. If a chord was classified correctly[7] we call it a true positives (TP) and visualize it in black. If a chord is classified incorrectly, this causes two errors. One error occurs at the annotated chord, which was not classified correctly. We call that a false negative (FN) and assign it the color red. Another error occurs at the classified chord which was not annotated. We call this error a false positive (FP) and display it in blue.

In general, we see that the system recognizes many chords correctly. However, there are also several errors. We can explain some of them musically, such as the error occurring between frames 40 and 50. Here, C♯m is annotated but EM is detected. Looking in the musical score we find that actually $\mathrm{C}^{\sharp}\mathrm{m}^{7}$ was played. In the annotation, the chord was mapped to C♯m but since $\mathrm{C}^{\sharp}\mathrm{m}^{7}$ comprises the pitch classes of both C♯m and EM, both results are possible. There are also other effects which can lead to a misclassification of a chord, such as overtones and the presence of non-chord notes. This kind of error is usually present for the entire duration of a chord or at

[7] In reality annotations are not always correct or unambiguous. Nevertheless we call a detected chord that matches the annotation "correct".

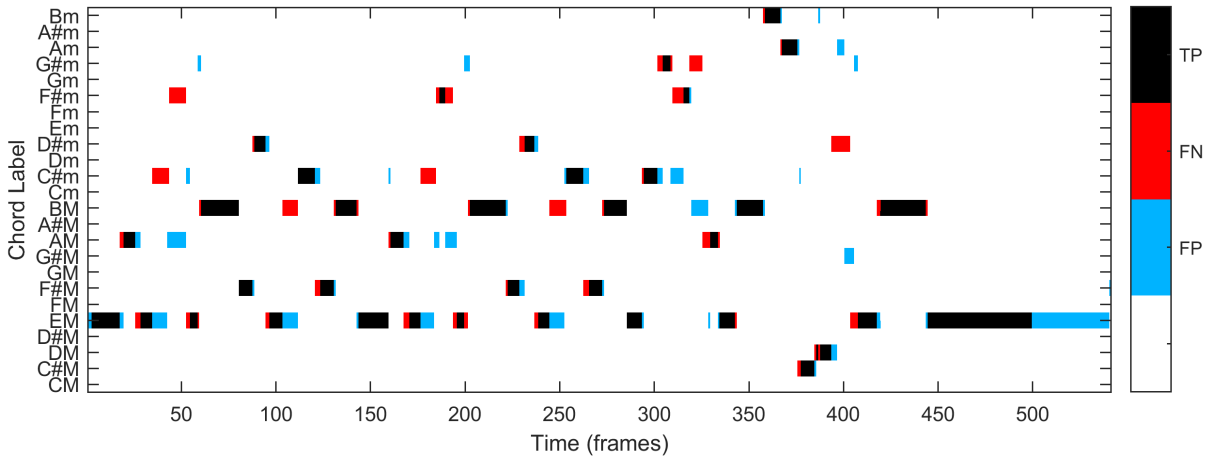


Figure 2.6: Results of template-based chord recognition of the Bach example with temporally smoothed chroma vectors.

least the duration of a note. However, there are many short errors lasting only a few frames. These errors result from the presence of noise in the signal. We can clearly recognize the errors since the duration of the detected chords is shorter than of naturally occurring chords, which we assume to be about one or two seconds. We can prevent that error by smoothing the chroma features over time before the classification. In Figure 2.6, a moving average with a two-sided window size of 11 frames was applied to the chromagram before classification. We see that most of the outliers are gone, but on the other hand, we now suffer from a reduced temporal resolution. There are more errors at the edges of correctly recognized chords. The overall accuracy was reduced from 75.9% without smoothing to 73.4% with smoothing. This smoothing approach does not respect the context of the recording. We introduce a more sophisticated smoothing technique in the next section.

The accuracy of a chord recognition result is defined as the percentage of correctly estimated annotated chord frames. This means that we ignore frames without annotated chord for the calculation. An example for such frames are the frames in the end of the Bach example ($n > 500$), where we perceive only silence. This is necessary due to the lack of the "no-chord" label in our chord model.

At this point, we briefly discuss typical chord confusions. The more common notes two chords have, the more likely the system will confuse them. Limiting ourselves to the major and minor chords, each chord has three chords which two common notes. For CM those are Cm, Em and Am. Note that the mode changes for all confusions. The root notes of the confusion chords are the original root note, a major third above the root note and a major sixth above the root note. We can find similar patterns for minor triads.

In this thesis we test several different chord vocabularies comprising more than just the 24 major and minor triads. This approach yields a problem if triads and tetrads are mixed. To illustrate that problem let us look at the defined similarity measure from a different perspective. With the templates defined in Section 2.1, the definition of the similarity measure is equivalent to averaging the energy of all pitch classes contained in a chord. Note that $\|\boldsymbol{x}\|_1 = 1$, by definition and $\|\mathbf{t}^\lambda\|_1$ equals the number of pitch classes contained in this chord. If we compare the similarity measures for CM and CM^7 we can also write:

$$\begin{aligned} s_\mathrm{t}\left(\boldsymbol{x}, \mathbf{t}^\mathrm{CM}\right) &= \frac{1}{3}\left(x_0 + x_4 + x_7\right) \\ s_\mathrm{t}\left(\boldsymbol{x}, \mathbf{t}^{\mathrm{CM}^7}\right) &= \frac{1}{4}\left(x_0 + x_4 + x_7 + x_{10}\right) \end{aligned} \tag{2.14}$$

Here, x_i denotes the i^{th} element of the chroma vector. These equations imply that CM^7 will only win if x_{10} is greater than the average of the remaining three elements, which is a very strict condition to be fulfilled and not a suitable criterion to decide whether a seventh chord is played. Note that we can alter the behavior by using different normalization schemes (e.g. no normalization, ℓ_2 norm, or element-wise weighting), but that would require another optimization procedure which might lead to over-fitting. Since there is no straightforward solution for this problem, we apply template matching only on vocabularies comprising chords with the same number of notes.

It is also possible to model a chord as probability distribution for chroma vectors. We can define a probability density function (PDF) $P(\boldsymbol{x}|\lambda)$ for each chord λ. A common choice for this PDF is a Gaussian function. This function depends on the mean vector $\boldsymbol{\mu}_\lambda$ and the covariance matrix Σ_λ, which have to be defined for each chord. There are basically two strategies to obtain those parameters. In [5], Bello and Pickens model the parameters using ideas similar to the ones we used for creating the templates with additional thoughts about the covariance matrix. This model is based only on musical knowledge. On the other hand in [9, 20] the authors learn the parameters from annotated data. [9] gives a more detailed analysis of this method including the usage of mixture models which are a superposition of several Gaussian PDFs.

When we use a Gaussian chord model, the score of a chord is not defined as a similarity measure, but rather as the value of the corresponding PDF evaluated at the current chroma vector. The further processing is equivalent to the template-based similarity measure. We compute the score for each chord and choose the chord with the highest score. In Figure 2.7, we show the results of the Gaussian-based chord recognition when applied to the Bach example. Here, we use two different recordings of the same choral to train the Gaussian models using a maximum likelihood estimation. The results are very similar to the results of template matching. The Gaussian model was able to match a few more chords and we reach an accuracy of 79.1%, which is slightly higher than the template based approach, but the problem of the small outliers remains. We

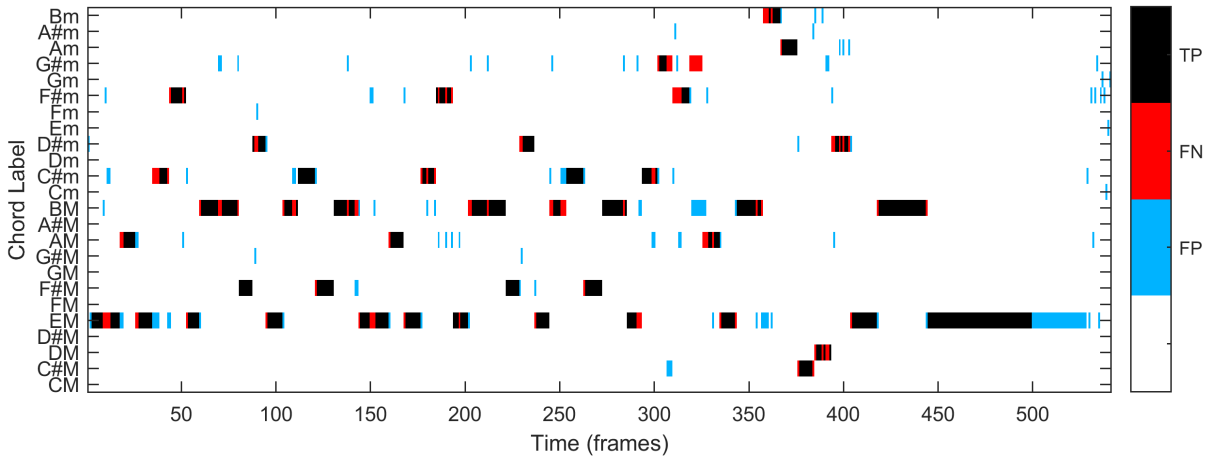


Figure 2.7: Results of the chord recognition with Gaussian chord models.

could not be solve the problem with this approach. The next section introduces a model that allows context-sensitive smoothing.

2.4 HMM-Based Chord Recognition

This chapter introduces the *Hidden Markov Model* (HMM), which is widely applied within mathematics and engineering [6, 10, 32] and can also be used for chord recognition [36]. An HMM is a model that describes sequence data by hidden internal states and an observable output variable. In the first part of this section, we explain the general idea behind the model and show that it can be useful for chord recognition. In the next part, we show the application to our problem. The theoretical part is mainly based on [33] and [24, Chapter 5.3].

The HMM is based on the so-called *Markov chain*. A Markov chain consists of a set $\mathcal{A}$ of I discrete states $\alpha_i \in \mathcal{A}$, $i \in [1:I]$. At each time step n, the model is in exactly one state $s_n \in \mathcal{A}$, $n \in [1:N]$, with N as length of the sequence. After each time step, the state changes according to a probability distribution which only depends on the previous state and neither on the states before the previous state nor on the specific time instant. This assumption is central in the theory of Markov chains and HMMs and is called the *Markov property*. We can write it as:

$$P(s_n = \alpha_j | s_{n-1} = \alpha_i, s_{n-2} = \alpha_k, \ldots) = P(s_n = \alpha_j | s_{n-1} = \alpha_i) = a_{ij}, \tag{2.15}$$

where the state transition probability $a_{ij} \in [0, 1]$ denotes the probability of a transition between α_i and α_j. We can summarize the state transition probabilities in the state transition matrix A. Together with the initial state probabilities $c_i = P(s_1 = \alpha_i)$, which describe the state probabilities at $n = 1$, we can compute the probabilities for all states at all times.

In many applications, the states of a system are not directly observable. Instead, we observe a variable that depends on the state in a probabilistic sense but is unable to tell us the state exactly. By hiding the states from the observer and introducing an observation variable $\boldsymbol{o}_n$, we arrive at the Hidden Markov Model. The observation variable $\boldsymbol{o}_n$ describes the observable output of the model at time n and follows a random distribution that depends only on the current state. We define $\mathcal{B}$ as the set of all possible observations and the state-dependent emission probability distribution $b_i(\boldsymbol{o}_n) : \mathcal{B} \to \mathbb{R}_{\geq 0}$, which maps the observation space $\mathcal{B}$ to the non-negative real numbers[8]. We define

$$b_i(\boldsymbol{o}_n) = P(\boldsymbol{o}_n | s_n = \alpha_i)\,, \tag{2.16}$$

where $\boldsymbol{x}$ is the observed vector. Overall, the HMM Θ is defined as a tuple containing the set of states $\mathcal{A}$, the transition matrix A, the set of initial state probabilities C, the set of observations $\mathcal{B}$ and the set of emission probabilities B:

$$\Theta = (\mathcal{A}, A, C, \mathcal{B}, B) \tag{2.17}$$

With this in mind, we now apply an HMM to chord recognition. We can use the chords as hidden states and chroma vectors as observations. We now show how the different aspects of an HMM are applicable on chord recognition. We first look at the chords, which we model as a Markov chain. We can split the recording in time frames and assign each frame one chord. The chord in the next time step depends on the previous chord. This is true since some chords are more likely to follow each other. For example transitions with fifth intervals between root notes are more probable than others. The most important transition to consider is the self-transition, i.e. the probability that the chord does not change. This probability is high compared to the other probabilities since a chord is usually played for more than one frame. On the other hand, the probability for a certain chord also depends on chords further in the past, since there are typical chord sequences in some genres, like the 12 bar blues scheme. However, such a behavior violates the Markov property and therefore exceeds the model. Therefore the HMM cannot take these properties into account. As for the initial state probabilities, we typically set them to a uniform distribution, since we usually have no prior knowledge about the scale or other tonal information about the recording.

We see that we can model the chords as a Markov chain. Since we do not observe the chord directly we extend the model to an HMM by adding the chroma vectors $\boldsymbol{x}$ as observations. If the HMM is in the state of a certain chord, this changes the probability distributions of the chroma vectors. If the model is in state CM for example, a chroma vector with high energies at C, E and G has a higher probability than other vectors. In Figure 2.8 we show a toy example of an HMM.

[8] For a finite number of observations, we can give actual probabilities, which only assumes values between 0 and 1. For continuous observations we need to use a probability density function, which can assume all non-negative real numbers.

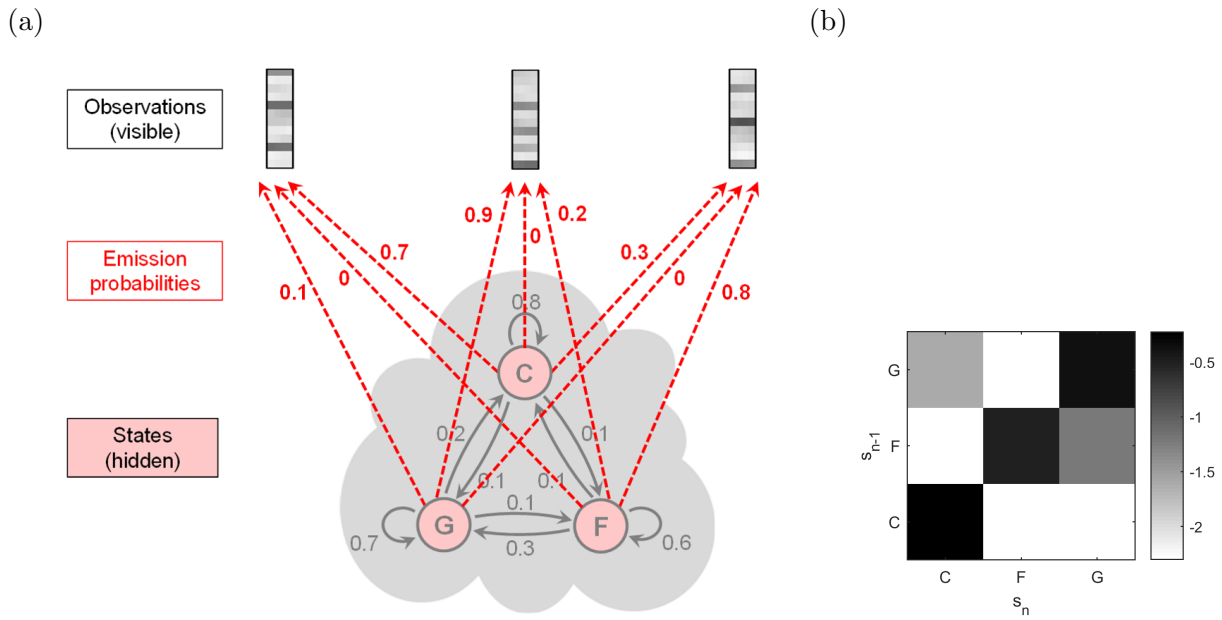


Figure 2.8: Toy example of an HMM describing only three chords. We show a state transition diagram with emissions in (a), taken from [24, Figure 5.28a]. We show the state transition matrix in (b). We plot the logarithm of the probabilities for a better visual quality.

The model comprises only three chords and three possible outputs. When we look at the state transition probabilities, we see that each chord favors a certain chroma vector output, but it is impossible to be sure about the chord just by looking at the chroma vector. The figure also shows a plot of the transition matrix in a way we use throughout the thesis. The vertical axis shows the previous state and the horizontal axis shows the target state.

With these considerations, we obtain a full hidden Markov model describing the chords in connection with the emissions. We can now reformulate the chord recognition problem as finding an optimal state sequence given a certain observation sequence. In fact there is a standard algorithm for solving this problem: the Viterbi algorithm. We explain the algorithm more thoroughly in Section 3.1. Figure 2.9 gives an example for an estimated chord sequence, which we computed with the Viterbi algorithm. Comparing the plot to Figures 2.5 and 2.6, there are less outliers but the edges of the chords are still preserved. The achieved accuracy rises to 83.3%. However, we still see the same errors as in Figure 2.5 at frames 40 to 50 and 170 to 200. Those can be explained musically and are hard to avoid due to ambiguities in the chord model.

Now we need to find suitable model parameters A and B. We can train the transition probabilities using labeled data. In [20], Jiang *et al.* counted the transitions in a training set to get an estimate. That way, they can exploit the general musical structure since many songs follow similar chord patterns. However, in [9], Cho and Bello showed that the main gain of HMM-based

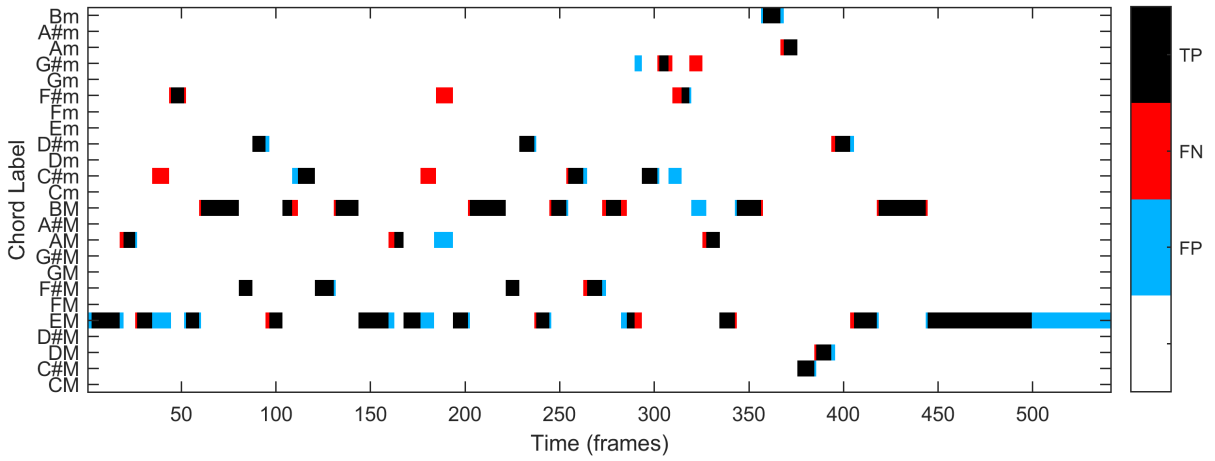


Figure 2.9: Chord sequence of the Bach example estimated with the Viterbi-algorithm.

chord recognition schemes is the smoothing property caused by a high self-transition probability. We can exploit the smoothing property by setting the main diagonal to a suitable value and setting all remaining entries to another, smaller value. In [5], Bello and Pickens propose another approach, which chooses the different transition probabilities according to the distance in the circle of fifths. The transition probabilities decrease linearly with higher distances. This captures some musical properties in a generic way without training. However, this method becomes more complicated when considering a larger number of chords, therefore in the following, we only enhance the main diagonal of the state transition probability matrix to achieve smoothing. Figure 2.10 shows two examples of possible transition matrix, one of the is trained from labeled data and the other one is modeled with an enhanced main diagonal. The trained matrix shows also higher probabilities for rising and falling fifth transitions which complies to music theory. The modeled matrix does not capture such properties.

We can use basic similarity measures as model for the emission probabilities, such as the template similarity measure s_t or the value of a Gaussian PDF modeling the chords. That way, we have two different models. In this thesis we call them "template emission model" and "Gaussian emission model". Note that both functions are not PDFs, since their integral over $\mathcal{F}$ is not equal to 1, as would be the case for an actual PDF. However, this is not important in the used algorithms. In the remainder of the thesis we limit ourselves to template emission models due to their simplicity and because they do not require training, but yield chord recognition results of sufficient quality.

This section introduced a powerful tool to jointly model chords and chroma vectors mathematically. With the HMM it is possible to obtain optimal chord sequences, using context-sensitive smoothing. We also can get insight in the probabilities for certain chords and chord transition, which we use in the next chapter.

(a) Trained Transition Matrix

(b) Modeled Transition Matrix

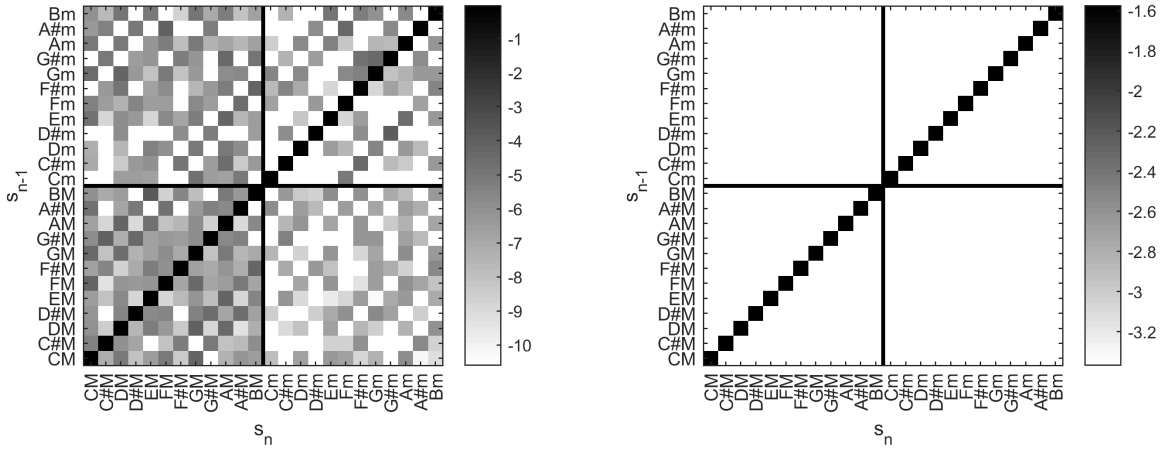


Figure 2.10: Two examples of possible transition matrices. We trained the matrix in (a) on annotated Beatles songs and modeled the matrix in (b) by enhancing the main diagonal of an uniform matrix. We display both matrices in the logarithmic domain.

Chapter 3

Chord Transition Features

In this chapter, we propose two different kind of chord transition features, which contain information about the frequencies of different chord transitions. In [45] Weiß proposes chord transition features and showed that information about chord transitions is useful for several classification tasks. The features need to be transposition invariant, i.e. invariant to recordings played in different keys.

In [45] Weiß estimated the chord sequence with the Vamp plugin "Chordino"[1], which uses the Viterbi algorithm, and counted the transitions. However, there was a hard decision for a chord sequence involved. Even if two chords have almost the same probability, the system has to decide for one before further processing. A wrong decision yields two errors, once in the transition into the chord, once in the transition to the next chord. In this thesis, we propose a feature extraction scheme that avoids hard decisions, aiming towards robust features with a high descriptive power. The features should be able to describe chord transitions in a way that depends only on the musical content and not on the specific recording, which may vary in tempo, key and instrumentation.

In the following, we propose two analogous transition feature extractors, which both are built on HMMs. The first one uses the Viterbi algorithm involving hard decisions, the other uses the Baum–Welch algorithm, which allows for a soft decision approach. The base for both approaches is a HMM modelling the chords as described in the Section 2.4.

3.1 Viterbi-Based Features

First, we introduce Viterbi-based chord transition features, which involve hard decisions. The *Viterbi algorithm* [43] is commonly used to determine the optimal state sequence in an HMM

[1] `http://www.isophonics.net/nnls-chroma`

Θ when the observation sequence is known. As discussed in the previous chapter, we can use this algorithm to estimate an optimal chord sequence given a sequence of chroma vectors. In this section, we introduce an approach which derives chord transition features from the detected sequence. At first, we introduce the Viterbi algorithm before discussing the feature generation itself. The following explanation is a summary of [24, 33].

The Viterbi algorithm finds an optimal state sequence $\hat{S} = (\hat{s}_1, \hat{s}_2, \ldots, \hat{s}_N)$, with N being the length of the sequence, also called *path*, such that the joint probability of the state sequence $S = (s_1, s_2, \ldots, s_N)$ and the observation sequence $O = (\boldsymbol{o}_1, \boldsymbol{o}_2, \ldots, \boldsymbol{o}_N)$ given a model Θ is maximized:

$$\hat{S} = \underset{S}{\operatorname{argmax}} P(O, S|\Theta)\,. \tag{3.1}$$

This is a hard optimization problem since testing all possible state sequences has a complexity of $\mathcal{O}\left(I^N\right)$, with I as the number of states. However, the Viterbi algorithm uses dynamic programming to decrease the complexity to $\mathcal{O}\left(I^2 N\right)$, which is linear in the sequence length N.

The idea behind the algorithm is to iteratively optimize the paths from the beginning of the sequence up to a certain state. This is done efficiently by using the results of the previous time step. This is only possible because of the Markov property, since otherwise we would need to look further back in the path. With the Markov property, however, we can optimize the path step by step by optimizing sub-sequences. If we add another state to the sub-sequence—which is already optimized—we only have to check all I^2 transitions from the last step to the next. We perform this procedure iteratively for all time steps $n \in [2 : N]$ until the whole sequence is optimized.

As an input, the Viterbi algorithm needs the observation sequence and the model parameters, particularly the number of states, the state transition matrix and the emission probability functions. Additionally, we need an auxiliary variable $D\,(i, n)$ which captures the probabilities of the optimized sub-sequences:

$$D\,(i, n) = \max_{s_1 \ldots s_{n-1}} P(\boldsymbol{o}_1 \ldots \boldsymbol{o}_n, s_1 \ldots s_{n-1}, s_n = \alpha_i|\Theta)\,. \tag{3.2}$$

$D\,(i, n)$ is the joint probability for the best (most probable) state sub-sequence ending in state α_i at time n and the given output sub-sequence $(\boldsymbol{o}_1, \boldsymbol{o}_2, \ldots, \boldsymbol{o}_n)$. Here, $n \in [1 : N] := \{1, 2, \ldots, N\}$ is the time index and $i \in [1 : I]$ is the state index. We can express the maximum path probability with the help of $D\,(i, N)$:

$$\max_{S=(s_1 \ldots s_N)} P(S, O|\Theta) = \max_{i \in [1:I]} D\,(i, N)\,. \tag{3.3}$$

We can compute $D\,(i, N)$ iteratively. The iteration starts by setting

$$D\,(i, 1) = P(\boldsymbol{o}_1, s_1 = \alpha_i|\Theta) = c_i b_i(\boldsymbol{o}_1). \tag{3.4}$$

To compute $D(i,n)$ for the following frame, consider that we know the maximal path probabilities to all states of the previous frame. As an intermediate step, we compute the path probability if for fixed states s_{n-1} and s_n:

$$\max_{s_1 \dots s_{n-2}} P(\boldsymbol{o}_1 \dots \boldsymbol{o}_n, s_1 \dots s_{n-2}, s_{n-1}{=}\alpha_j, s_n{=}\alpha_i|\Theta) = D(j,n-1) \cdot a_{ji} \cdot b_i(\boldsymbol{o}_n), \tag{3.5}$$

with $j \in [1:I]$. We compute the probability by multiplying the maximum probability to reach α_j in frame $n-1$ (given by $D(j,n-1)$) with the transition probability from state α_j to α_i (given by a_{ji}) and the probability that state α_i emits $\boldsymbol{o}_n$ (given by $b_i(\boldsymbol{o}_n)$). We can derive the expression using Bayesian inference. By maximizing this expression over s_{n-1} we obtain the following formula for $D(i,n)$:

$$D(i,n) = b_i(\boldsymbol{o}_n) \cdot \max_{j \in [1:I]} D(j,n-1) \cdot a_{ji}. \tag{3.6}$$

With this iteration procedure we can compute $D(i,N)$. However, this does not yet solve the problem since we search the sequence itself, not its probability. Therefore, we need to retrace our steps to arrive at $D(i,N)$. Algorithmically, we implement the procedure by introducing a backtracking variable $E(i,n) \in [1:I]$, which saves the states yielding the optimal sequences corresponding to each time step n and state s_i. We define it as

$$E(i,n-1) = \operatorname*{argmax}_{j \in [1:I]} \left(D(j,n-1) \cdot a_{ji} \right). \tag{3.7}$$

To apply the backtracking, we define state indices $i_n \in [1:I]$, such that the sequence $S = \alpha_{i_1} \alpha_{i_2} \dots \alpha_{i_N}$ is optimal. Starting with

$$i_N = \operatorname*{argmax}_{j \in [1:I]} D(j,N), \tag{3.8}$$

we apply recursion to find the remaining indices:

$$i_n = E(i_{n+1}, n). \tag{3.9}$$

That way we can find the entire optimal sequence:

$$\hat{s}_n = \alpha_{i_n}. \tag{3.10}$$

We can visualize one central step of the Viterbi algorithm using a so-called *trellis diagram* [2] in Figure 3.1. A trellis diagram shows the states for each time step as compared to a state transition diagram as in Figure 2.8. The trellis diagram is useful to visualize paths and time-dependent variables such as $D(i,n)$. We illustrate the idea of the Viterbi-algorithm by showing a toy example with three states. As an example we compute $D(3,n)$. To calculate $D(3,n)$ the paths

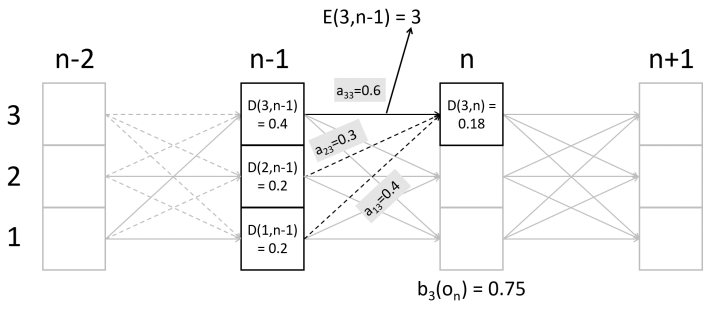


Figure 3.1: Iteration step of the Viterbi algorithm.

from all states in the previous time step are considered. The path with the highest score—the "survivor" [43]—which we illustrate by a solid line is used to calculate $D(3, n)$. All other paths are ignored (dashed lines). That way the path is always optimal until time step n. All survivors are possible parts of the final optimal sequence. The state from which the survivor originates is saved in the backtracking variable $E(i, n-1)$. The dashed lines between the time steps $n-2$ and $n-1$ indicate that there is only one allowed path to each state in time step $n-1$. Using such a diagram, we can obtain the optimal path by retracing the solid lines starting at the state with the highest probability. Computationally, we use the backtracking variable. Note that not all survivors are reached, since there are I survivors per time step.

As described in Chapter 2.4, we can now use the Viterbi algorithm to estimate an optimal chord sequence. For this, we use the chord labels as states and the chroma vectors as observations. After we obtained an optimal chord sequence as output we visualize the detected chords by a slightly modified trellis diagram in Figure 3.2. We color those states that are part of the optimal sequence in black, the other states are white. We can now visualize the transitions by connecting temporally adjacent states of the optimal sequence. We draw all transitions as red lines. We see a very stable behavior. The majority of transitions are self-transitions.

We now obtained knowledge about the local chord transitions. We use those features to generate piece-level features in Chapter 3.4.

3.2 Baum–Welch-Based Features

In the previous chapter, we derived local features describing the chords and chord transitions based on the Viterbi algorithm. The Viterbi algorithm returns exactly one chord label for each time step—the information about all other states is discarded. We use the sequence to derive local chord transition features by "connecting" temporally adjacent chords. In this section, we introduce another kind of feature which avoids hard decisions. Instead of deriving the

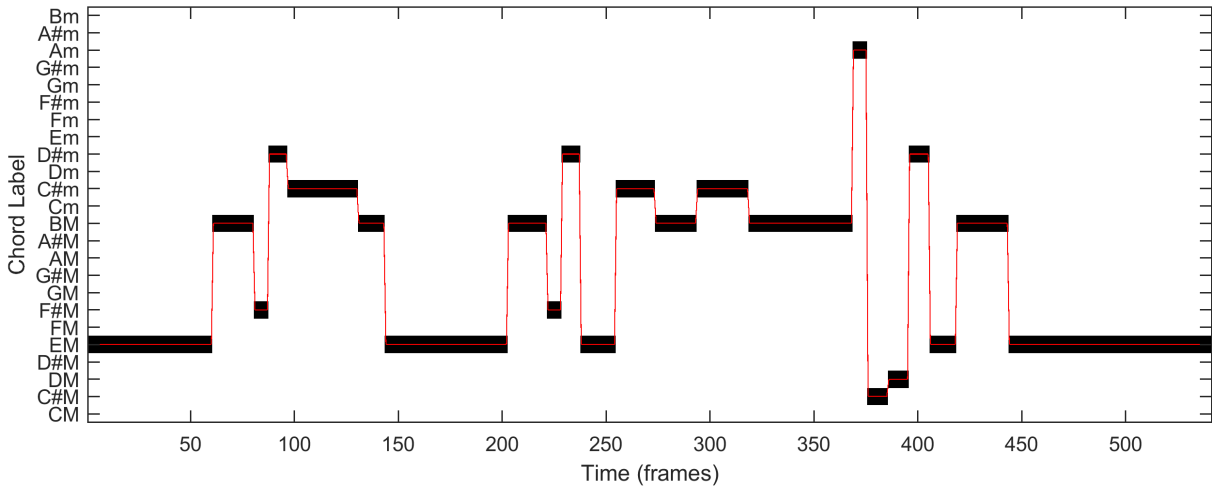


Figure 3.2: Visualization of the Viterbi-based chord transitions using the Bach example.

transitions from the most probable chord sequence, we compute probabilities for all transitions in each time step. This information is contained in an intermediate variable of the *Baum–Welch algorithm*, which was originally developed for training HMMs [3]. We first describe the Baum–Welch algorithm already in the context of chord transition estimation. Afterwards, we show visualization techniques for the obtained local chord transition features.

3.2.1 Baum–Welch Algorithm

Baum proposed the algorithm [3], which is now known as Baum–Welch algorithm, as a method to re-estimate model parameters in HMMs. The following explanation of the Baum–Welch algorithm is based on the tutorial of Rabiner [33]. We start with an initial model $\Theta = (\mathcal{A}, A, C, \mathcal{B}, B)$, specifying all parameters. Furthermore, the algorithm requires an observation sequence O. The algorithm estimates a new model $\hat{\Theta} = \left(\mathcal{A}, \hat{A}, \hat{C}, \mathcal{B}, \hat{B}\right)$, which explains the observation sequence better than the old model, such that

$$P\left(O|\hat{\Theta}\right) \geq P(O|\Theta)\,. \tag{3.11}$$

The parameters are estimated from $\xi_n\,(i,j)$ which contains the probabilities that a transition from α_i to α_j occurs at time step n. This variable is our major interest in the Baum–Welch algorithm, since it can directly serve as local chord transition feature. The transition probability is computed using the probabilistic models of HMMs. Formally, we define:

$$\xi_n\,(i,j) = P(s_n = \alpha_i, s_{n+1} = \alpha_j | O, \Theta)\,, \tag{3.12}$$

	$v_n(i)$	$w_n(i)$
Initialization	$v_1(i) = c_i b_i(\boldsymbol{o}_i)$	$w_N(i) = 1$
Iteration	$v_{n+1}(j) = \left[\sum_{i=1}^{I} v_n(i)\, a_{ij}\right] b_j(\boldsymbol{o}_{n+1})$	$w_n(i) = \sum_{j=1}^{I} a_{ij} b_j(\boldsymbol{o}_{n+1}) w_{n+1}(j)$

Table 3.1: Equations of the forward and backward algorithm which are used to compute $v_n(i)$ and $w_n(i)$ respectively.

for $n \in [1 : N-1]$. To compute the probability for a transition, we need to consider the probability that we arrive in state $s_n = \alpha_i$ together with the previous observations. This probability is captured in the forward variable $v_n(i)$:

$$v_n(i) = P(\boldsymbol{o}_1 \boldsymbol{o}_2 \ldots \boldsymbol{o}_n, s_n = \alpha_i | \Theta)\,. \tag{3.13}$$

Furthermore, we need the probability that we observe the future observations given that we are in state $s_{n+1} = \alpha_j$. We define backward variable $w_n(i)$

$$w_n(i) = P(\boldsymbol{o}_{n+1} \boldsymbol{o}_{n+2} \ldots \boldsymbol{o}_N | s_n = \alpha_i, \Theta)\,, \tag{3.14}$$

which represents this probability.

In order to compute $v_n(i)$ directly, we have to sum up the joint emission probabilities of all possible paths ending in $s_n = \alpha_i$, which is computationally not feasible. Similarly, to compute $w_n(i)$, we have to sum up the emission probabilities of all paths starting in $s_n = \alpha_i$. Using dynamic programming, we can decrease the computational effort. There are iterative algorithms, the forward and backward algorithm, to compute both variables efficiently. We state them in Table 3.1 without further proof. We derive the equations for both algorithms in Appendix A.1 and A.2. Here, we give an intuitive explanation of the forward algorithm. For a more detailed discussion of both algorithms, please refer to [33].

The forward algorithm exploits the Markov property. Using this property it is computationally feasible to sum over all possible paths if we know the forward variables of the previous time step. In turn, those variables contain information about all paths leading to this time step and so on. Figure 3.3 shows our toy example, this time we compute one step in the forward algorithm. We see that we have to consider only three different transitions for the computation of $v_n(3)$. The information about the history of the sequence is already contained in the forward variables of the previous time step. We see, that for capturing the many different paths leading to $s_n = \alpha_3$, only the information drawn in black is needed. This visualizes the low computational complexity. The intuition behind the forward algorithm is very similar to the Viterbi algorithm. Also the formulas are analogous, the maximization over all states is simply replaced by a sum over all

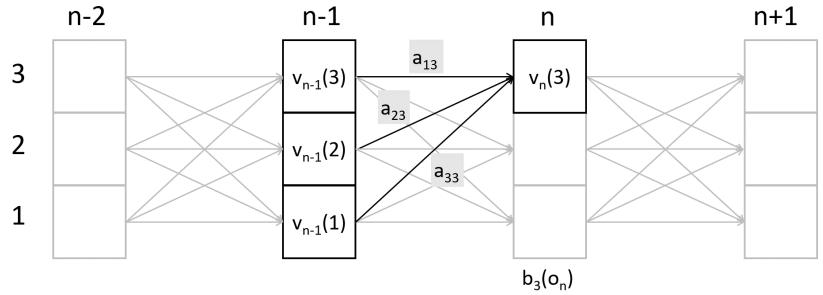


Figure 3.3: This figure shows the calculation of $v_n(3)$ out of the forward variables from time step $n-1$. All relevant information is drawn black whereas the rest of the trellis diagram is grey.

states. (Compare the iteration step of the forward algorithm with Eq. 3.6.) The considerations for the backward algorithm are very similar, with a reversed order.

We compute the forward and backward variables by initializing $v_1(i)$ and $w_N(i)$ for all $i \in [1:I]$ according to the equations in Table 3.1. Afterwards, we apply the iteration step for all $n \in [1:N]$ and $i \in [1:I]$. From those two variables, we can then compute the desired transition probabilities:

$$\xi_n(i,j) = \frac{v_n(i)\, a_{ij} b_j(\boldsymbol{o}_{n+1}) w_{n+1}(j)}{P(O|\Theta)}. \tag{3.15}$$

We derive this formula in Appendix A.3 using Bayesian inference and the Markov property. We can explain the expression for $\xi_n(i,j)$ intuitively, by considering the contributing components. The probability for a transition depends on the probabilities of all paths leading to state $s_n = \alpha_i$, captured by $v_n(i)$, as well as the probabilities of all paths leading away from state $s_{n+1} = \alpha_j$ which are represented by $P(\boldsymbol{o}_{n+1} \dots \boldsymbol{o}_N, s_{n+1} = \alpha_j|\Theta) = b_j(\boldsymbol{o}_{n+1}) w_{n+1}(j)$. Also, the prior probability for a transition from α_i to α_j, represented by a_{ij} must be taken into account. The denominator $P(O|\Theta)$ is merely a normalization factor and is needed for a probabilistic interpretation.[2] We can compute it by:

$$P(O|\Theta) = \sum_{i=1}^{I} v_N(i). \tag{3.16}$$

In Figure 3.4 we visualize the components needed to calculate $\xi_n(i,j)$. Here, we compute the transition probability of a single transition in our toy example. To calculate the probability $\xi_n(3,3)$ of the transition marked in red, all paths to $s_n = \alpha_3$ have to be considered, as well as all paths leading away from $s_{n+1} = \alpha_3$. Because of the Markov property, it is sufficient to consider all paths only up to one time step away. Those path segments are marked in black and are considered in the forward and backward variable respectively.

[2]The probability $P(O|\Theta)$ is the solution of the so-called evaluation problem as described in [33] and gives the probability to observe an observation sequence given a model. Using it we can determine which model out of several probably emitted an observation sequence. Solving the evaluation problem is useful for example, in speech recognition.

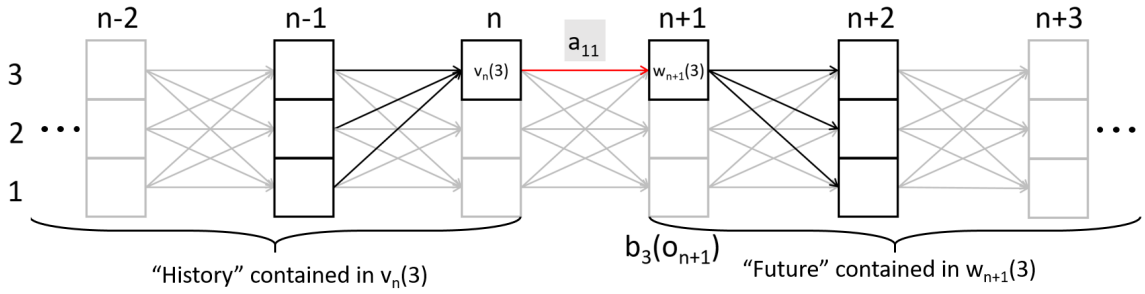


Figure 3.4: Schematic illustration of the calculation of $\xi_n(i,j)$, marking all relevant components in black.

If we know the probabilities of all transitions originating from α_i at time n, we can obtain the probability that any transition originates from α_i at time n, by summing up the probabilities. This probability is equivalent to the probability of being in state α_i at time n. The probability can be computed from $\xi_n(i,j)$ by marginalization:

$$\gamma_n(i) = P(s_n{=}\alpha_i|O,\Theta) = \sum_{j=1}^{I} \xi_n(i,j) = \frac{v_n(i)\, w_n(i)}{P(O|\Theta)}. \tag{3.17}$$

We can express $\gamma_n(i)$ also just in terms of the forward and backward probabilities. This equation can be derived from Eq. 3.15 and the iteration step of the backward algorithm[3]. Even though $\xi_n(i,j)$ is only defined for $n \in [1:N-1]$, this expression holds for all $n \in [1:N]$. At this point, we can stop the algorithm since we have information about the state probabilities and the state transition probabilities. We can use those probabilities to derive soft chord transition features and perform a kind of soft chord recognition. We summarize the procedure in Algorithm 3.1. Here, we use a data structures in form of multi-dimensional arrays $V(i,n)$, $W(i,n)$, $\Gamma(i,n)$, and $\Xi(i,j,n)$ instead of the time series variables $v_n(i)$, $w_n(i)$, $\gamma_n(i)$, and $\xi_n(i,j)$, respectively in order to assimilate MATLAB code.

For the sake of completeness we briefly discuss the remaining Baum–Welch algorithm. With $\gamma_n(i)$ and $\xi_n(i,j)$ we can re-estimate the HMM model parameters. By summing up $\xi_n(i,j)$ and $\gamma_n(i)$ over time, we obtain the expected number of transitions between two states and the expected number of occurrences of each state. We can re-estimate the new transition probabilities $\hat{a}_{ij}$ by relating both measures.

$$\hat{a}_{ij} = \frac{\text{expected number of transitions from } \alpha_i \text{ to } \alpha_j}{\text{expected number of transitions from } \alpha_i} = \frac{\sum_{n=1}^{N-1} \xi_n(i,j)}{\sum_{n=1}^{N-1} \gamma_n(i)}. \tag{3.18}$$

[3]This part of the Baum–Welch algorithm is actually called forward backward algorithm. Since it is strongly connected to the features derived from the Baum–Welch algorithm and we use it to gain information about them we still refer to this as Baum–Welch algorithm to avoid confusion.

Algorithm: Baum–Welch-based chord transition estimation
Input: HMM specified by $\Theta = (\mathcal{A}, A, C, \mathcal{B}, B)$
Observation sequence $O = (\boldsymbol{o}_1, \boldsymbol{o}_2, \ldots, \boldsymbol{o}_N)$
Output: Local state probabilities $\Gamma(i, n)$
Local state transition probabilities $\Xi(i, j, n)$

Procedure:
Forward Algorithm: Initialize an $(I \times N)$ 2D-array V by $V(i, 1) = c_i b_i(\boldsymbol{o}_1)$ for $i \in [1 : I]$. Then, compute in a nested loop for $n = 1, \ldots, N-1$ and $j = 1, \ldots, I$:

$$V(j, n+1) = \left[\sum_{i=1}^{I} V(i, n)\, a_{ij}\right] b_j(\boldsymbol{o}_{n+1})$$

Set the normalization factor: $\nu = \left[\sum_{i=1}^{I} V(i, N)\right]^{-1}$

Backward Algorithm: Initialize an $(I \times N)$ 2D-array W by $W(i, N) = 1$ for $i \in [1 : I]$. Then, compute in a nested loop for $n = N-1, \ldots, 1$ and $i = 1, \ldots, I$:

$$W(i, n) = \sum_{j=1}^{I} a_{ij} b_j(\boldsymbol{o}_{n+1}) W(j, n+1)$$

Initialize an $(I \times N)$ 2D-array Γ and set

$$\Gamma(i, n) = \nu \cdot V(i, n)\, W(i, n)$$

Initialize an $(I \times I \times N)$ 3D-array Ξ and set

$$\Xi(i, j, n) = \nu \cdot V(i, n)\, a_{ij} b_j(\boldsymbol{o}_{n+1}) W(j, n+1)$$

Return: $\Gamma(i, n)$ and $\Xi(i, j, n)$

Algorithm 3.1: Baum–Welch-based chord transition estimation

We can find new estimations for the initial state probabilities $\hat{c}_i$ by:

$$\begin{aligned} \hat{c}_i &= \text{expected number of times in state } \alpha_i \text{ at time n} = \\ &= \gamma_1(i)\,. \end{aligned} \tag{3.19}$$

The re-estimation of the emission probability functions depends on the used probabilistic model. For a Gaussian model, we can re-estimate mean vector and covariance matrix as follows:

$$\hat{\boldsymbol{\mu}}_i = \frac{\sum_{n=1}^{N} \gamma_n(i)\, \boldsymbol{o}_n}{\sum_{n=1}^{N} \gamma_n(i)} \tag{3.20}$$

$$\hat{\Sigma}_i = \frac{\sum_{n=1}^{N} \gamma_n(i)\, (\boldsymbol{o}_n - \boldsymbol{\mu}_i)\, (\boldsymbol{o}_n - \boldsymbol{\mu}_i)^{\intercal}}{\sum_{n=1}^{N} \gamma_n(i)}, \tag{3.21}$$

which is a straightforward way of estimating Gaussian models. In HMM training applications, we initialize the parameters either randomly or by an educated guess. Afterwards, we perform the algorithm iteratively, each time with an updated model until the model converges. In this

thesis, however, we do not use the Baum–Welch algorithm for training, which is why we do not cover the training step in detail. Here, we use the algorithm for soft-decision chord recognition and the corresponding transition estimation. We refer to [33] for further information about the Baum–Welch algorithm and its application on HMM training.

3.2.2 Visualization Techniques

To use the Baum–Welch algorithm for chord estimation, we initialize the HMM in the same way as for the Viterbi-algorithm. Afterwards, we use the algorithm to compute local state probabilities $\gamma_n(i)$ and state transition probabilities $\xi_n(i,j)$. We can visualize these probabilities similar to Figure 3.2 by connecting all states with red lines representing the respective transitions. Instead of coloring only one state per time frame in black, we now color all states with a gray value indicating the probability of being in that state. A darker tone represents a higher probability. Also, since there is now more than one possible transition at each time step, we have to plot the transition according to the probabilities as well. In a first approach, we plot every transition with a probability exceeding a certain threshold. The transparency of the line indicates the probability of the respective transition. A more intense red color refers to a high transition probability. In order to properly show the chord transitions, it is necessary to ignore the self-transitions, since they are much more probable and would visually suppress the other transitions. We show the result in Figure 3.5a. We see that the plot is not easily readable. This is because the differences between the values are too small to be captured by such a color map. There are many chords with similar, high probabilities, e.g. between EM and C♯m, which share two common notes. In this case we see that the transitions EM → C♯m and Em → C♯M are strong in all frames where both chords are present. This is an artifact of the Baum–Welch algorithm and cannot be avoided. The transitions going back and forth between states create a symmetric, lattice-like structure. While those transitions may have musical reasons, they do not describe relevant harmonic change in the recording. On occasions where the harmonic content changes, we observe more directed transitions, which are visually covered by the noise-like transitions.

In oder to visualize the harmonic changes, we need to find an algorithm which enhances the musically relevant, directed transitions, and suppresses the lattice-like "noise floor". We have three types of transitions: The self-transitions, which have the highest probabilities, the noise-like transitions, which have the lowest probabilities, and the musically relevant probabilities. We assume that those transitions stand out by slightly higher probabilities. Filtering out the probabilities that stand out from the noise is a complicated problem for multiple reasons. A major issue is that we deal with normalized matrices. Because of the normalization, the value of each element is dependent on all other elements. As a result, the probability ranges of the small value of ρtioned transitions types vary from frame to frame. We also do not know how many transitions we want to enhance since there might be more than one transition of interest.

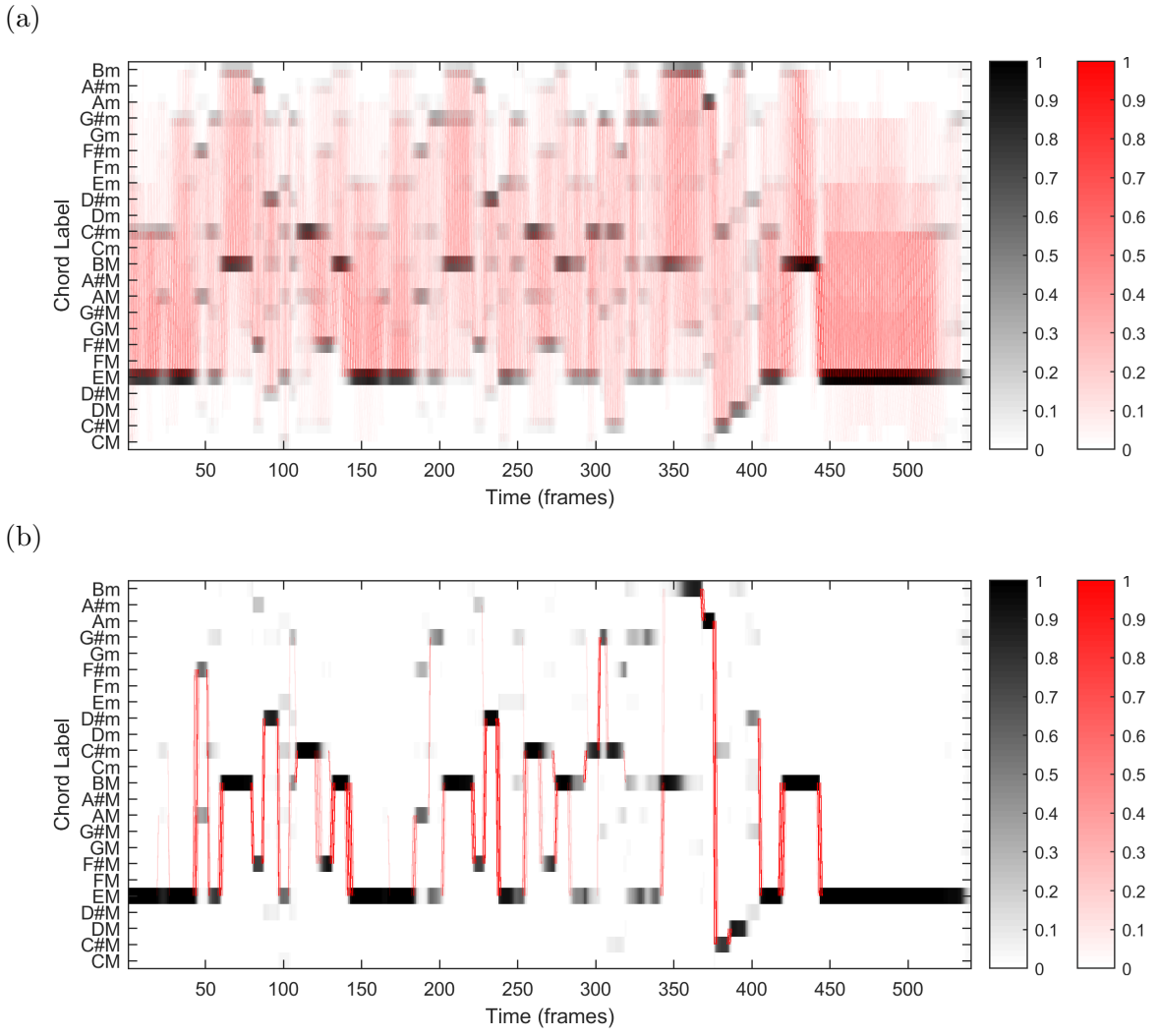


Figure 3.5: Straightforward (a) and enhanced (b) plot of the state and transition probabilities.

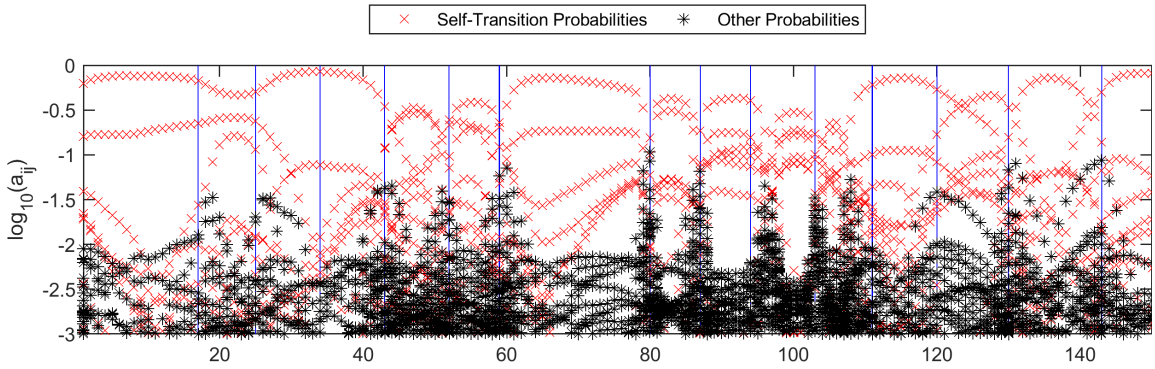
Since we know which matrix elements describe the self-transitions, we can discard them, but distinguishing the other types of transitions is harder. It is not sufficient to raise the threshold, since the scales differ between the time frames. A way to circumvent that problem is the so-called softmax function. This function enhances the higher values of a given set while suppressing the lower ones. We define the softmax function for a matrix $X \in \mathbb{R}^{M \times N}$ as follows:

$$\mathrm{softmax}_\eta\left(X\right) = \frac{\exp\left(\eta X\right)}{\sum_{m=1}^{M}\sum_{n=1}^{N}\exp\left(\eta X\left(m,n\right)\right)}, \tag{3.22}$$

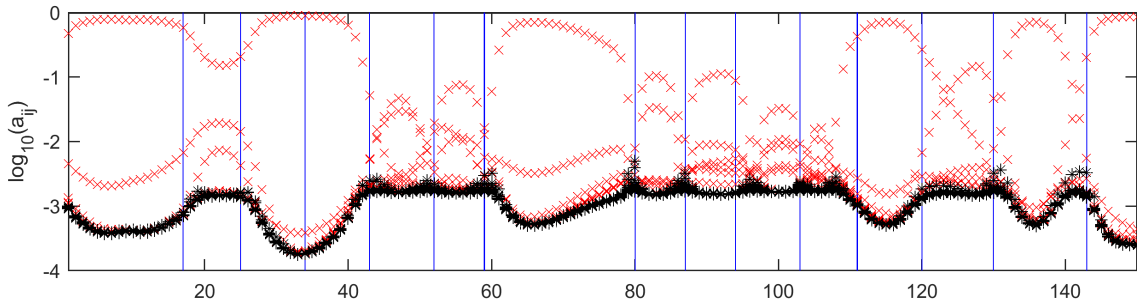
where $\exp(\cdot)$ is the element-wise exponential function and $\eta \in \mathbb{R}_{>0}$ is a parameter determining the "softness" of the function. A smaller η corresponds to a softer maximization. If we apply the softmax function on the transition probabilities of each time frame—after deleting the self-transition probabilities—we enhance the highest probabilities in a flexible way. However, if there is no value standing out from the noise, the function enhances the highest value, which is part of the noise.

At this point we see that the complexity of the rescaling task is too high for a conventional approach like a single softmax rescaling. We therefore introduce an empirically found algorithm: the double-softmax rescaling. As the name indicates, the algorithm employs two consecutive softmax functions. At first, we apply the softmax function with a small η_1 on all transition probabilities of one frame, including the self-transition probabilities. Since there is usually at least one high self-transition probability, this leads to a compression of the remaining probabilities. Afterwards, we set all self-transition probabilities to zero, before we apply another softmax function with a high η_2 on the resulting matrix. This enhances the highest of the remaining off-diagonal probabilities. We demonstrate the functionality of the double-softmax rescaling in Figure 3.6. For this figure we plot all entries of the transition probability matrix for each time frame in the logarithmic domain. That way, we can see the distribution of the probabilities for each frame. We indicate the positions of the annotated chord changes by vertical blue lines. In the first figure, we plot the distribution as obtained by the Baum-Welch algorithm. We color the self-transition probabilities in red and the other transition probabilities in black. We see that we have one high self-transition probability in each frame, which corresponds to the chord with the highest probability. When we look at the other transitions we see a "noise floor" with several outliers. However, there are many outliers—also in frames without annotated chord transitions—and we cannot be sure how many of them are musically relevant. In the second plot, we show the values after the first softmax function. We see that all off-diagonal transition probabilities are strongly compressed and follow a line that is inverse to the sequence of highest self-transition probabilities. I.e. if there is one high self-transition probability, the values of the other transitions are smaller. This is a typical behavior of normalization within the softmax function. We now take a closer look at the other transitions in the third plot by zooming in the y-axis. We observe outliers only in a few frames now and also the number of the outliers

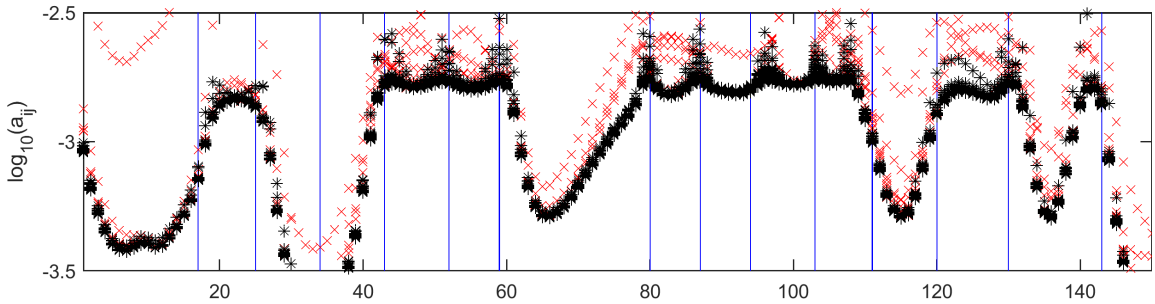
(a) No Rescaling



(b) After the First Softmax



(c) After the First Softmax (Zoomed)



(d) After the Second Softmax

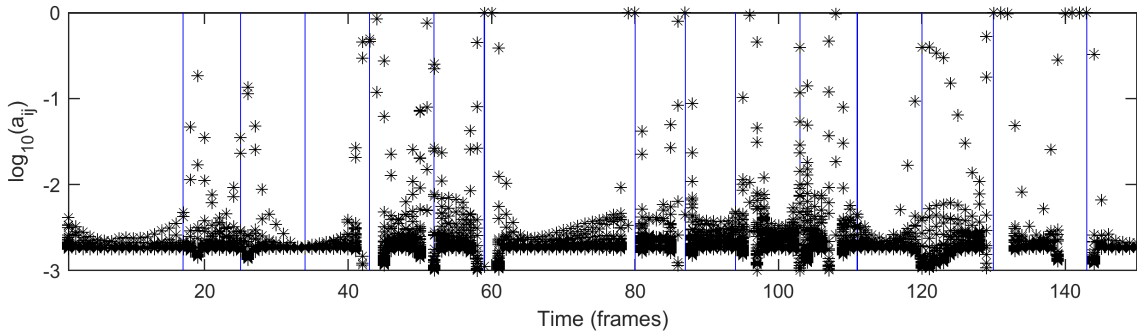


Figure 3.6: Distribution of the transition probabilities for each frame. The blue lines indicate annotated chord transitions.

Algorithm: Double-Softmax Rescaling
Input: Temporal sequence of $N-1$ state transition matrices $\xi_n(i,j)$
Softmax parameters η_1, η_2
Output: Temporal sequence of $N-1$ state transition matrices $\xi_n(i,j)$ with suppressed lattice structure
Procedure:
In a loop over all time steps $n \in [1:N-1]$:
Initialize a $(I \times I)$ matrix Ξ with $\Xi(i,j) = \xi_n(i,j)$
Set $\Xi = \text{softmax}_{\eta_1}(\Xi)$
Set $\Xi(i,i) = 0$ for all $i \in [1:I]$
Set $\Xi = \text{softmax}_{\eta_2}(\Xi)$
Replace $\xi_n(i,j)$ with $\Xi(i,j)$ for all $i,j \in [1:I]$

Return: $\xi_n(i,j)$

Algorithm 3.2: Double-softmax rescaling algorithm for the suppression of the lattice structure.

is considerably smaller than before. We see that the outliers match the annotated transitions to some extent. The values in most frames are now strongly suppressed. The second softmax enhances those outliers, such that only a few strong transitions remain. If the values are too close to each other or too small, the second softmax does not enhance these transitions. We show the resulting distribution after the whole procedure in the fourth plot. We clearly see the noise floor on the bottom and only a few transitions which have a value higher than 0.01. We can find those transitions mostly around the positions of the annotated chord changes.

We now plot those values instead of the original probabilities in our trellis diagram in Figure 3.5b. We empirically chose the parameters $\eta_1 = 10$ and $\eta_2 = 10000$. Additionally, we apply a single softmax ($\eta = 10$) on all state probabilities of each time frame. We see that we succeed in suppressing the lattice structure and that the remaining transition are in fact the transitions which connect the most probable states. In areas where the state is uncertain, we see transitions to both states. A good example is around frame 420 where we see transitions from EM to both BM and Bm. On the other hand, some transitions are not visible anymore since the state changes only gradually. As an example we look at the frames around time step 350, where the detected chord changes slowly from BM to Bm. Here, the transition BM $\rightarrow$ Bm was spread over several frames, so that the individual transitions are not strong enough to be enhanced. However, even though the double-softmax procedure yields good visual results, the tuning of the softmax parameters is difficult and depends on the used HMM parameters and the data.

In summary, the enhancement algorithm consists of three steps, which we subsequently apply on all individual matrices $\xi_n(i,j)$ for each time step $n \in [1:N]$. For a compact explanation, see the box Algorithm 3.2.

3.3 Comparison and Evaluation

In the following, we compare the Viterbi-based and the Baum–Welch-based local chord transition features. We do not have a quality measure for local transition features available, so we compare the transitions qualitatively. However, since both feature types are based on a chord recognition scheme using HMMs, we can quantitatively compare the performance of both algorithms in a chord recognition scenario.

3.3.1 Influence of the Self-Transition Probability

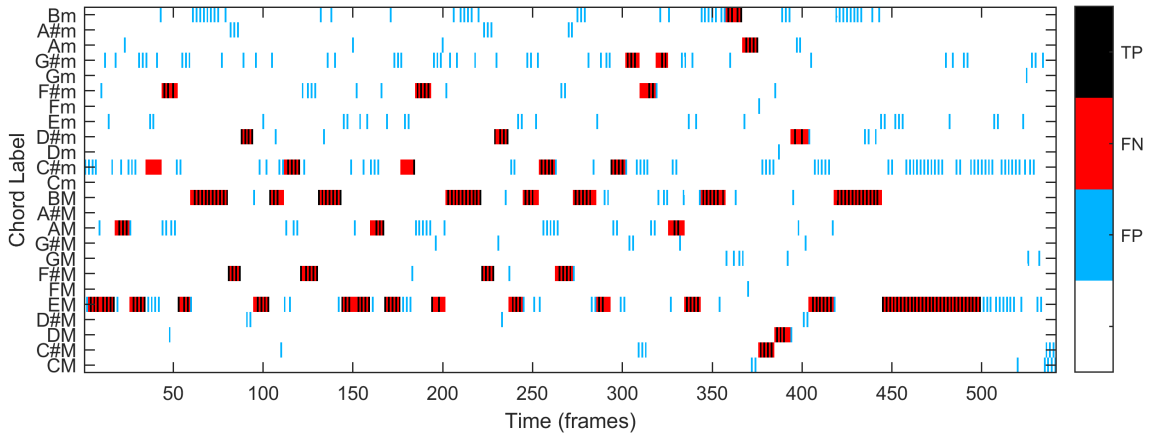
As a first experiment, we analyze the HMM parameters. In Chapter 2.4, we introduce the state transition matrix A. We show that we can model the matrix with a uniform matrix of which we enhance the main diagonal. Such a matrix depends on only one parameter, which describes the ratio between the main diagonal elements and the off diagonal elements. In [9], Cho and Bello defined such a parameter. However, we found that this definition has disadvantages when we compare models with different numbers of states since the self-transition probability depends on the number of states in the HMM. We propose a different parameter ρ which directly defines the self-transition probability. We assume that the smoothing properties stay the same even if we use chord vocabularies of different sizes and, therefore, self-transitions matrices with different sizes. In [33], Rabiner states that the average duration of a state in a Markov chain depends only on the self-transition probability of that state. Since the smoothing property is connected with the average state duration, this result supports our parameter definition. We define the state transition probabilities as follows:

$$a_{ij} = \begin{cases} \rho & \text{if } \; i = j \\ \frac{1-\rho}{I-1} & \text{if } \; i \neq j \end{cases}, \tag{3.23}$$

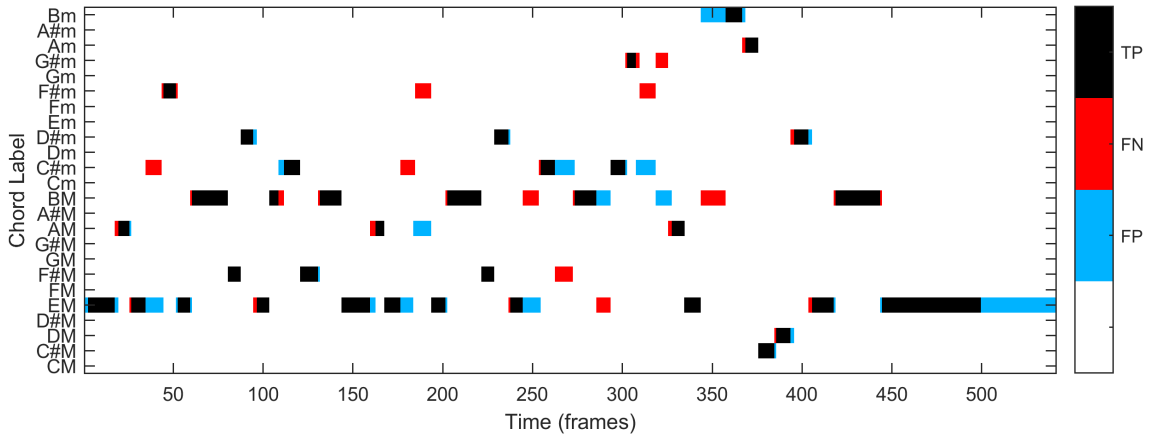
with the size of the chord vocabulary I and $0 \leq \rho \leq 1$. In the context of chord recognition, a higher choice of ρ leads to a longer average chord duration. Therefore, we obtain a more stable result. To get a rough idea of the influence, we show the results of the chord recognition for three selected values of $\rho \in \{0.025, 0.2, 0.5\}$ in Figure 3.7. In this experiment we use the Viterbi algorithm. For $\rho = 0.025$, the chord changes very often, we speak of over-segmentation. Apparently, the smoothing is not strong enough. Nevertheless, for $\rho = 0.5$, the algorithm misses several transitions and the smoothing is too strong. This is called under-segmentation. The value of $\rho = 0.2$ yields the best of the shown results. We show similar plots for the Baum–Welch algorithm in Appendix B.1.

We see that ρ has to be tuned carefully, since it influences the results greatly. Therefore we systematically examine the influence of ρ on a larger database. For this purpose, we use the

(a) $\rho = 0.025$



(b) $\rho = 0.2$



(c) $\rho = 0.5$

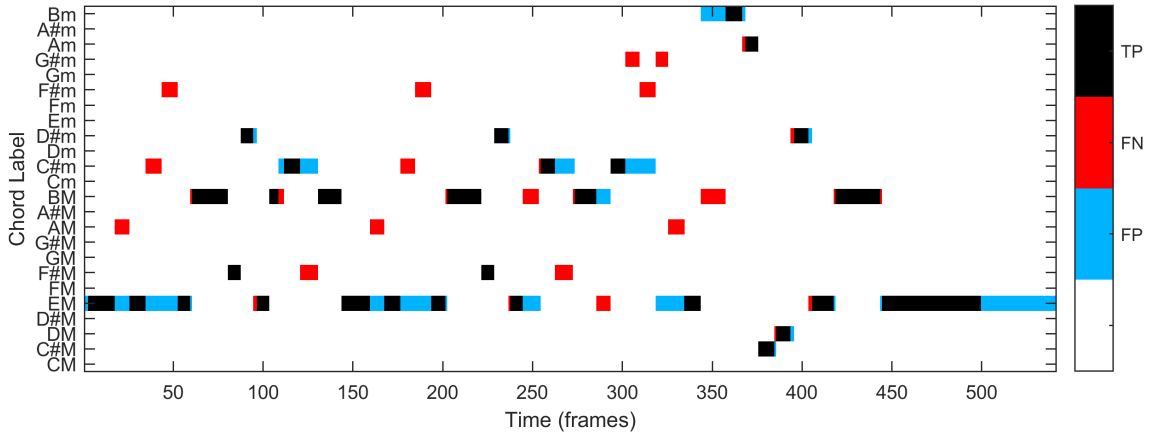


Figure 3.7: Viterbi-based chord recognition of the Bach example. We use different ρ to model the chord transition probabilities.

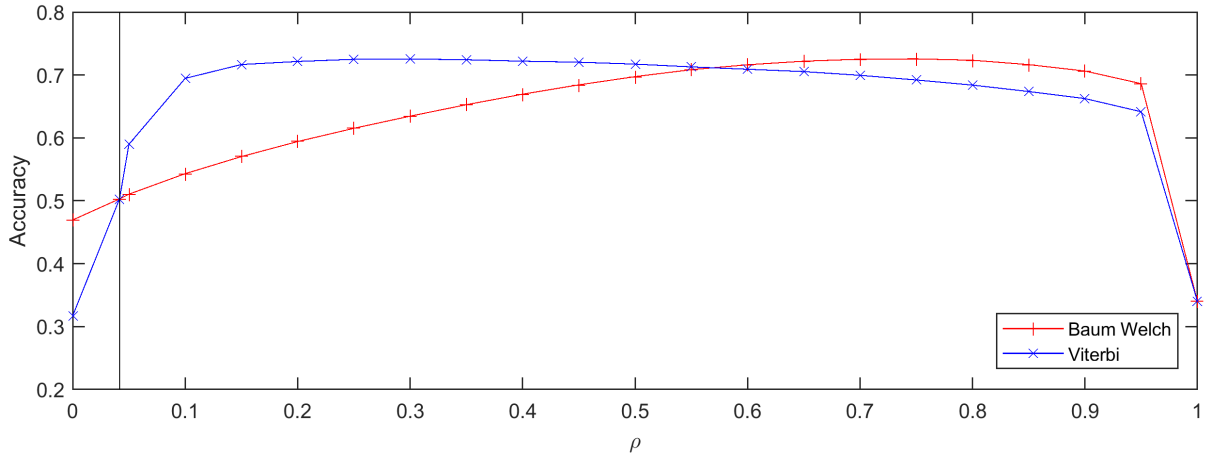


Figure 3.8: The accuracy of the chord recognition as function of ρ. The experiment was performed on the *Beatles* dataset using either the Baum–Welch algorithm or the Viterbi algorithm. The vertical line at $\rho = \frac{1}{24}$ marks the point where all transitions have equal probability.

Beatles dataset. It comprises 180 recordings of Beatles songs from the 10 studio albums. Detailed chord annotations for all songs are publicly available [22]. The dataset is widely used for benchmarking chord recognition systems within the MIR community [20]. To compare the results, we perform HMM-based chord recognition on all songs and compute the accuracy by comparing the result with the annotated ground truth. We define the accuracy as the ratio of correctly classified frames. Since our chord model does not contain a "no-chord" symbol, we discard all frames which were annotated with "no-chord". For the experiment, we reduce all annotated chords to the chord vocabulary of major and minor triads. We map each chord with a major third as first interval to the respective major triad and each chord that starts with a minor third to the respective minor triad. This is not always justified musically, especially when we deal with augmented and diminished chords, but is common practice [20]. The accuracy is then averaged over all songs. The result is used as a quality measure for the tested chord recognition system. For the evaluation of the Viterbi algorithm, we can compare the output sequence directly with the annotation. The Baum–Welch algorithm does not generate such a sequence. Instead, we use $\gamma_n(i)$ and choose the chord for which $\gamma_n(i)$ is maximal in each time frame

$$\hat{s}_n = \underset{s_n}{\operatorname{argmax}} P(s_n, O|\Theta) = \underset{\alpha_i}{\operatorname{argmax}} \gamma_n(i) \,. \tag{3.24}$$

We perform the described experiment with different values of ρ. Figure 3.8 shows the average chord recognition accuracy over all songs as a function of ρ. We see that the two algorithms show different behaviors. In particular, the optimal value of ρ for the Baum–Welch algorithm is higher than the optimal value of ρ value of the Viterbi algorithm. Since the Baum–Welch algorithm does not optimize the path probability but rather the individual state probability,

the algorithm allows other paths to enhance the probabilities of single states and is therefore less "strict" than the Viterbi algorithm, which optimizes the entire sequence. As a result, to obtain the same smoothing effect, we have to choose a higher self-transition probability for the Baum–Welch algorithm, which explains the different locations of the maximums.

However, we have to be careful when interpreting this result. The chord annotation is not unique, in particular with respect to the segmentation. In [24, Figure 5.16], Müller gives an example, how the song "Let it be" can be annotated by different music experts. If we fit the self-transition probability to an annotation which is highly segmented, the resulting optimal value of ρ will probably be smaller than a value of ρ fitted to an annotation with coarse segmentation.

Another noteworthy observation is the saturation of the accuracy for the Viterbi algorithm. For $\rho < 0.025$ the accuracy does not depend on ρ anymore. In this region, the self-transition probability is smaller than the other transitions, it is therefore "cheaper" to switch the state rather than to stay in the state. The Viterbi algorithm is therefore forced to change the state often in order to optimize the joint probability of the sequence. At a certain point, the chord changes in (nearly) every time step since staying in a chord is less probable then switching back and forth to the second most probable chord. Further decreasing the self-transition probability does not change that behavior, since no self-transitions exist anymore. Also, in the case of very small ρ, the Baum–Welch algorithm allows other paths to enhance the probability of a state. For a demonstration of this behavior, see Appendix B.1

The vertical line in Figure 3.8 marks the point where $\rho = \frac{1}{24}$ and thus, all transition probabilities are equal. The output only depends on the emission probabilities, since the model now does not inherently prefer one path to any other. In other words, we perform template matching. Therefore, the result does not depend on the algorithm anymore.

In general, the results from Figure 3.8 are consistent with the intuition we gained from Figure 3.7. There exists an optimal value for ρ, which is used as a default value for the future experiments. We set $\rho = 0.3$ for the Viterbi algorithm and $\rho = 0.75$ for the Baum–Welch algorithm. We refer to those values as "standard ρ". We see, that in these maximums the average accuracy is almost equal. We can conclude that in the scenario of chord recognition, none of the algorithms is better than the other. However, we see that the Viterbi algorithm is less sensitive on ρ and has a lower complexity than the Baum–Welch algorithm. Therefore, if we need only a discrete sequence of chords, the Viterbi algorithm should be preferred. However, if we are interested in a soft-decision output, we need to use the Baum–Welch algorithm.

We now use the obtained parameters to examine the Bach choral with both algorithms. We need to keep in mind that the parameter ρ was fitted on pop music, while we apply it on classical music. Since the two styles are very different the parameter might not be suitable. We show the result in Figure 3.9. There is not much difference when we look at the detected chords. Also, the

accuracy is similar with 71.2% for the Viterbi algorithm and 74.8% for Baum–Welch algorithm. The plot of the Baum–Welch result shows the probability that a chord is played at each time step. We use a softmax with $\eta = 10$ on the state probabilities to visually enhance the plot. While we see some time frames where the probability is almost one (e.g. frames 400–450), there are also several frames where two or more chords have similar probabilities. This is the case around time frame 50. There is a high probability both for F♯m and AM. This means that the algorithm was not sure about the played chord. In the score we see that actually the minor seventh chord F♯m^7 was played. In this special scenario, it is therefore possible to gain knowledge exceeding the model, since F♯m^7, which is not part of the used chord vocabulary, comprises pitch classes of both detected chords. This way, we are able to detect some seventh chords types. The Viterbi algorithm only detects F♯m, which is actually correct in the sense of the annotation, without giving a hint about the seventh chord.

3.3.2 Context-Sensitive Smoothing

The reason why we used HMMs in the first place was the context-sensitive smoothing. We argued that we can achieve a result where outliers are smoothed but sharp edges could be preserved at clear chord boundaries. To validate this theory, we perform a smoothed chord recognition in two ways: once by smoothing the chroma vectors with a moving average filter (window size of 21 frames) and a subsequent template matching (compare Section 2.3) and once with the Baum–Welch algorithm ($\rho = 0.97$). We choose the parameters empirically to obtain strong but similar smoothing. In both cases, the smoothing is much stronger than needed, but it helps us to observe the general smoothing behavior of both approaches. Figure 3.10 compares both algorithms. To get a more detailed view on the transitions, we plot only the first 150 frames. We show the frame-wise chord probabilities. In the first case we apply a softmax ($\eta = 150$) on the state probabilities to visualize the differences better. We see that the amount of smoothing is too high in both cases, but the Baum–Welch algorithm manages to obtain sharp edges and to precisely locate the chord transitions. An example can be found around frame 60. Even though the previous chord was smoothed too much, the transition to the next chord is very sudden and at the correct time. On the other hand, a simple smoothing of the chroma vectors yields a very smooth and inaccurate transition. Around frame 130 we see that the Baum–Welch algorithm classifies the chord falsely by extending the previous chord too long. However we still obtain a sharp edge on the transition to the next chord. This scenario visualizes the context-sensitive smoothing we can achieve with HMMs.

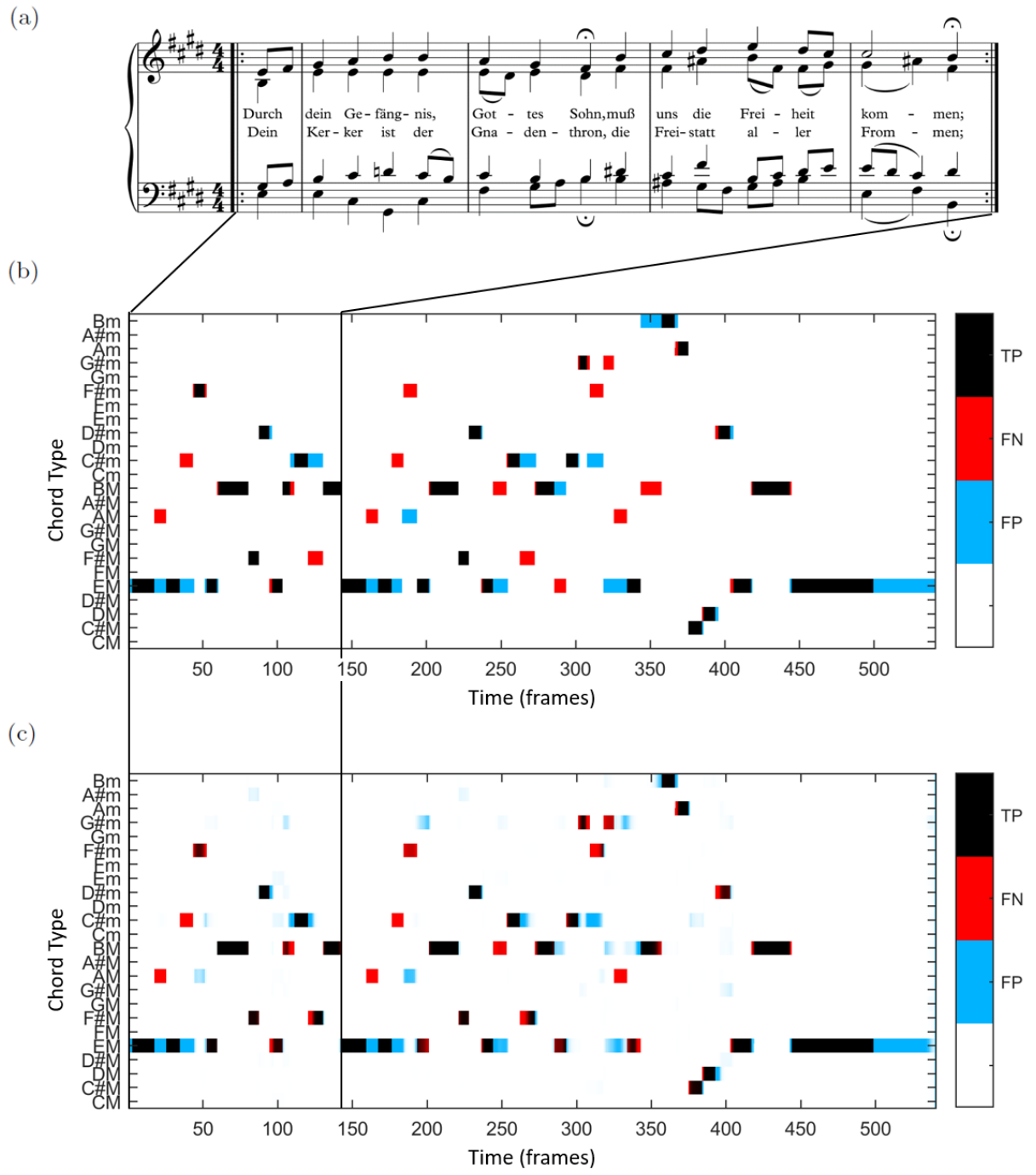


Figure 3.9: Comparison of the recognized chords when using (b) the Viterbi and (c) the Baum–Welch algorithm with standard ρ. The score of the fist part of the choral is shown in (a).

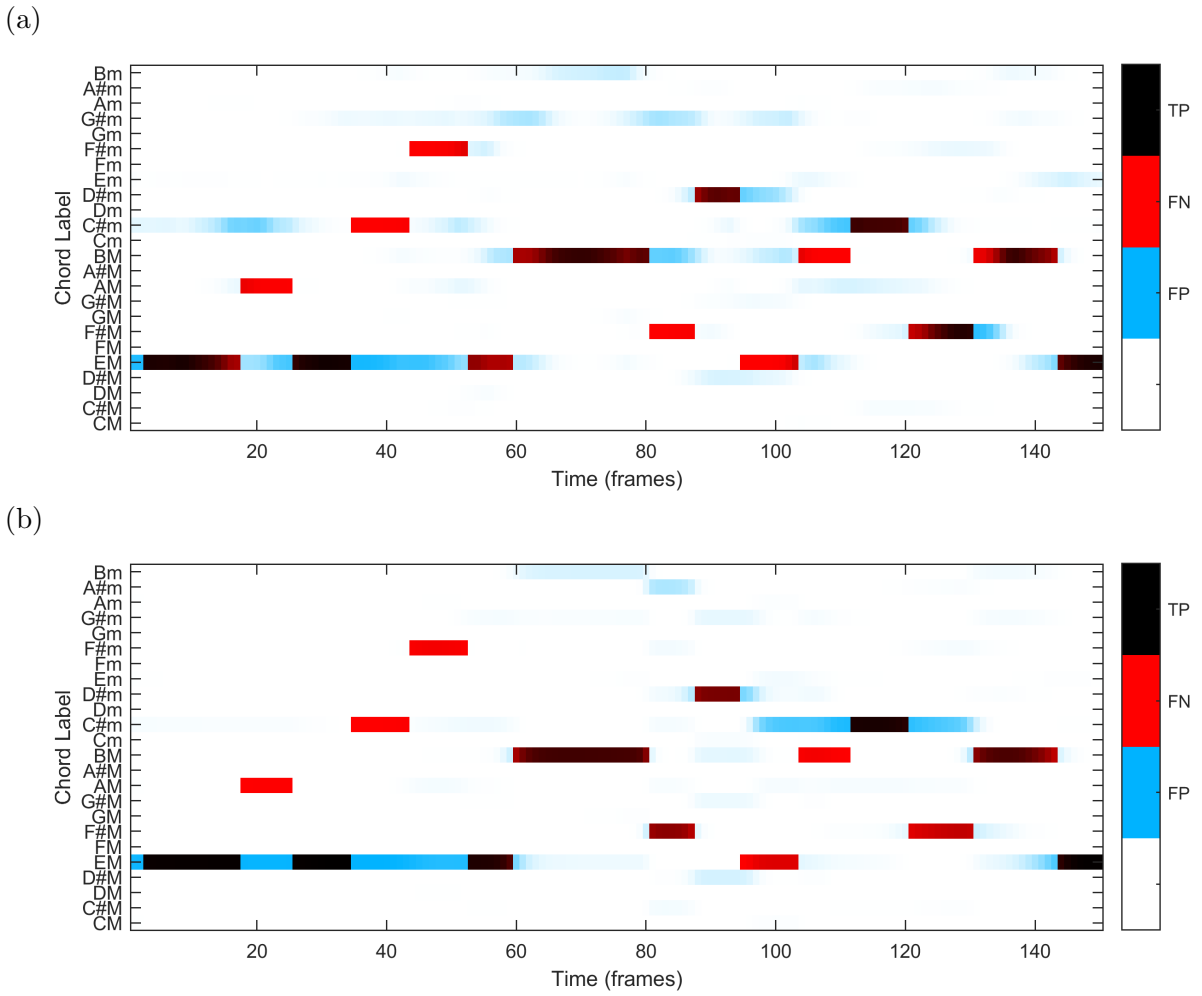


Figure 3.10: Chord recognition of the Bach example with strong smoothing. As smoothing algorithm we use (a) a moving average of the chroma vectors and (b) the Baum–Welch algorithm.

3.3.3 Local Chord Transition Features

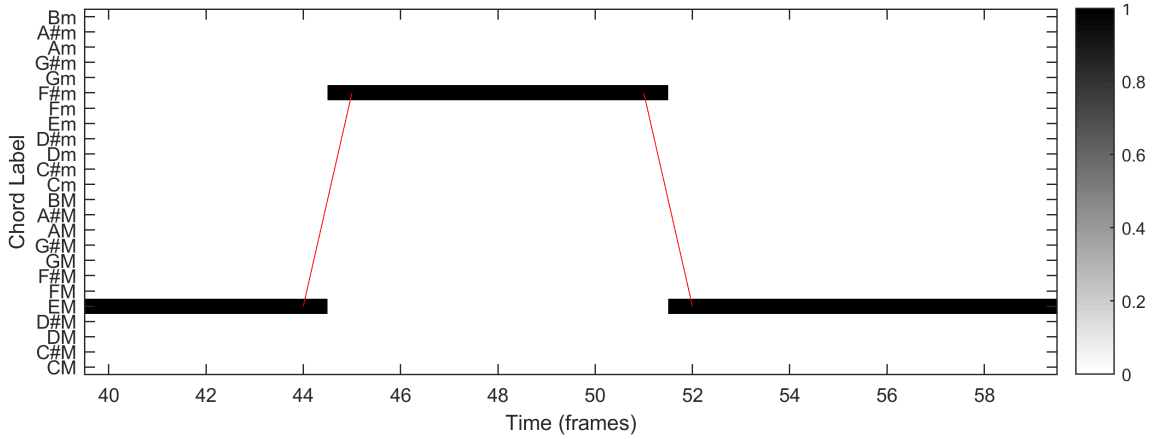
As a next step, we analyze the chord transitions computed with the Viterbi and the Baum–Welch algorithm. In particular, we look at the transitions around the small value of ρtioned seventh chord. In Figure 3.11, we show the chord transitions for the frames 40–59. The Viterbi algorithm yields only two transitions, whereas the Baum–Welch algorithm shows several. Not all of them describe harmonic changes. First of all, we see the expected transitions EM $\rightarrow$ AM and EM $\rightarrow$ F$^\sharp$m (frame 40–46). We see transitions in the other direction when the chord changes back to EM (frame 50–55). The transitions are spread over several frames. This indicates, that a transition can occur at each of those frames. We see, that the probability of the previous chord decreases as the probability of the next chord increases. At the same time, we see transitions from EM $\rightarrow$ G$^\sharp$m and back again. This is due to a confusion between EM and G$^\sharp$m, which share two notes. The same is true for Em and C$^\sharp$m. There are also several other chord transitions visible. It is not immediately clear why the corresponding chords occur. We can see those transitions as an artifact. The strength of the transitions contains some information about overtones and therefore relates to timbre. We see that we were not able to completely eliminate timbral content. On the other hand, the confusions allow for conclusions about the tonal content. For example, if only a fifth interval is played, this is usually mapped to the respective major chord due to the overtone structure. However, since no third is present, the confusion chords (and the corresponding transitions) except the major-minor confusion probably have only low probabilities. We can argue that those transitions do not have a clear semantic meaning anymore.

Using the double softmax rescaling introduced in the previous section we can suppress many transitions as shown in Figure 3.11(c). Here, only the transitions which relate to harmonic changes remain. We see the transitions from EM to both possible chords AM and F$^\sharp$m, and back. The plot only shows transitions which correspond to harmonic change but is still respecting that there might be more than one possible chord.

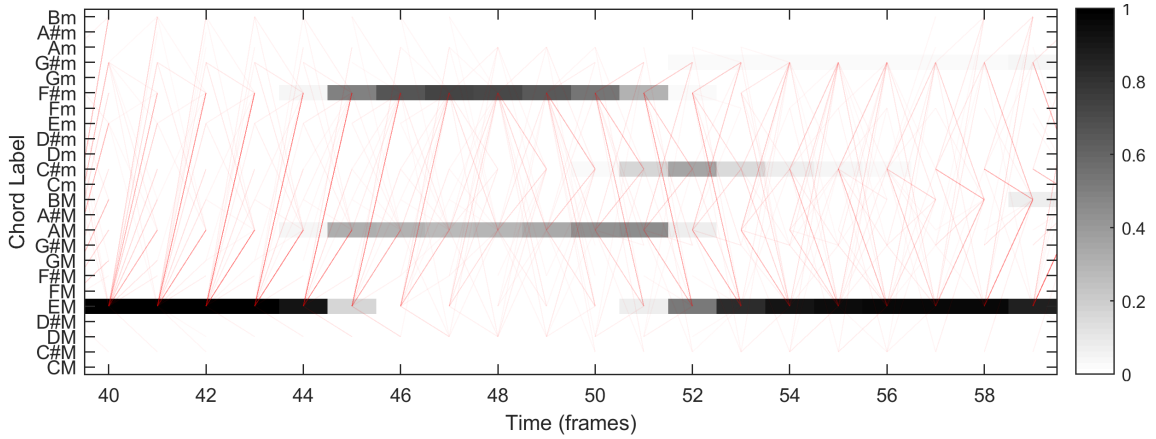
3.4 Piece-Level Chord Transition Features

In the previous sections, we discuss two different ways of extracting local information about chord transitions. The Viterbi algorithm yields an optimal state sequence from which we derive the local transitions. The Baum–Welch algorithm produces transition probabilities for each time frame.

(a) Viterbi



(b) Baum–Welch



(c) Baum–Welch (Double-Softmax)

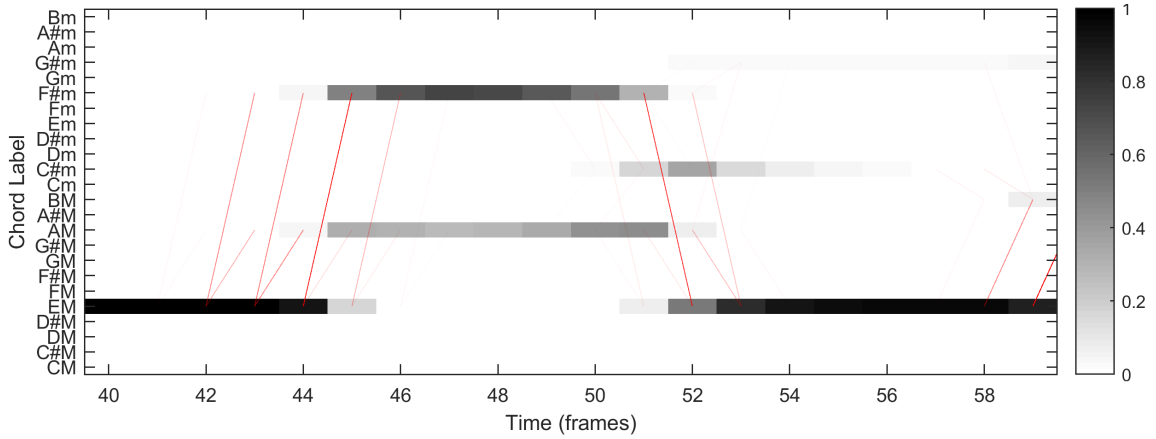


Figure 3.11: Comparison of the chord transitions when using (a) the Viterbi and (b) the Baum–Welch algorithm with standard ρ. In (c) we show chord transitions computed with the Baum–Welch algorithm after a double-softmax rescaling.

3.4.1 Feature Aggregation

In this section, we explain the aggregation algorithm which we use to aggregate the local chord transition features into global piece-level features, which describe a whole recording instead of single time instances. At first, we are interested in the relative frequency of occurrence of each chord transition. We define the variable a_{ij}^{w} that describes the average probability of a transition from λ_i to λ_j:

$$a_{ij}^{\mathrm{w}} = \frac{1}{N-1} \sum_{n=1}^{N-1} P(s_n = \lambda_i, s_{n+1} = \lambda_j)\,. \tag{3.25}$$

We call a_{ij}^{w} weighted chord transition probabilities. The name is based on the state transition probabilities a_{ij} defined for an HMM. We can interpret a_{ij}^{w} as a kind of state transition probability[4] which is weighted with the probability of occurrence of the previous chord. As a consequence, we have

$$\sum_{i=1}^{I} \sum_{j=1}^{I} a_{ij}^{\mathrm{w}} = 1, \tag{3.26}$$

as opposed to,

$$\sum_{i=1}^{I} a_{ij} = 1 \quad \forall j \in [1:I]. \tag{3.27}$$

The main difference is that a_{ij} is a conditional probability and a_{ij}^{w} is a joint probability. We denote the matrix comprising all weighted chord transition probabilities as A^{w}.

Starting from the optimal chord sequence obtained from the Viterbi algorithm, we can count the transitions and set:

$$a_{ij}^{\mathrm{w}} = \frac{\text{Number of transitions from } \alpha_i \text{ to } \alpha_j}{\text{Total number of transitions}}. \tag{3.28}$$

We need to use a different approach when we use the Baum–Welch algorithm. We obtain the Baum–Welch-based chord transition probabilities a_{ij}^{w} by averaging the local transition probabilities over all $N-1$ pairs of neighboring time steps. We can compute the chord transition matrix directly by:

$$A_{ij}^{\mathrm{w}} = \frac{1}{N-1} \sum_{n=1}^{N-1} \xi_n\,(i,j)\,. \tag{3.29}$$

From this point on, the procedure does not depend on the chord transition estimation algorithm anymore. A^{w} contains piece-level information about relative chord transition frequencies, which we could use as a feature already. However, it is not robust enough against different versions, since it contains more information than needed. As we mention in the introduction of this chapter, the feature should be invariant under key and tempo changes. Therefore, we discard all information about those two measures. The chord transition matrix contains implicit information

[4] For the Viterbi-based features, this is actually a relative frequency rather than a probability.

about the key of the recording, since certain chords are preferred in certain keys. For example, a recording in the key of C major most likely contains the CM triad more often than the C♯M triad. This is also reflected in the chord transition probabilities. We eliminate the influence of the key by performing a suitable transform to make the transitions transposition invariant (TPI), similar as shown in [24, Chapter 5.3.4.2]. We call this procedure the TPI transform. In the end, a transition from a chord should be independent from the root note of that chord. Since we deal with different chord types, we have to treat transitions between chord types separately. To achieve this, the matrix is split into sub-matrices such that each sub-matrix contains only transitions from one chord type to another chord type. In the case of only major and minor triads, this results in four sub-matrices (major → major, major → minor, minor → major and minor → minor). For more chord types, the number of sub-matrices increases accordingly. Also, note that—since there are only four distinct augmented chords—those sub-matrices may also be of non-square shape. We transform the sub-matrices separately by averaging over all cyclically shifted versions. Let X be an arbitrary sub-matrix. Then we can write:

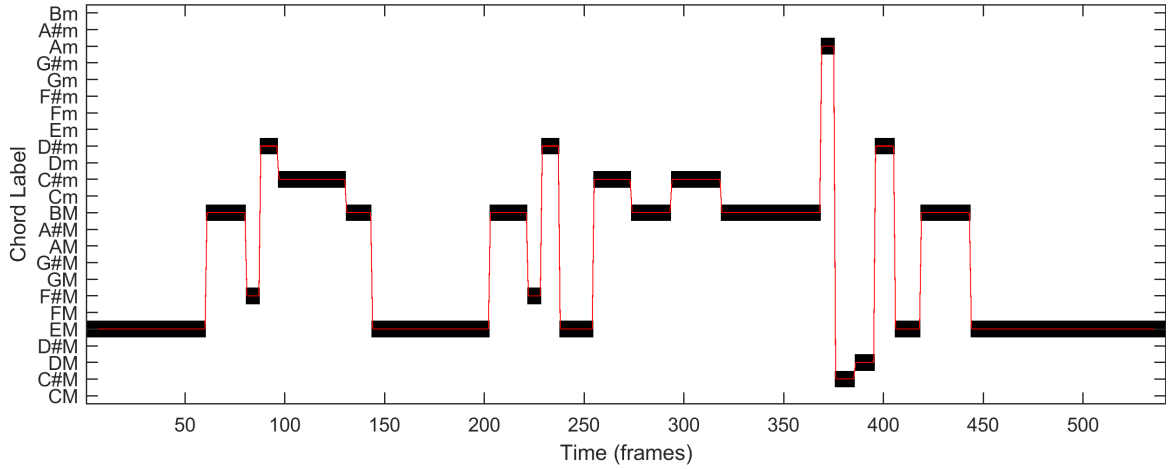
$$X_{ij}^{\mathrm{TPI}} = \frac{1}{12} \sum_{l=0}^{11} X_{(i+l) \bmod 12,(j+l) \bmod 12}. \tag{3.30}$$

Afterwards, we rebuild the matrix A^{TPI} from the sub-matrices X^{TPI}. That way an entry in the matrix no longer refers to a transition between two specific chords. Rather, each entry describes a transition specified by an interval between root notes and a change of chord type. For example, $a_{1,5}^{\mathrm{w}}$ is the probability of a transition from CM to EM, whereas $a_{1,5}^{\mathrm{TPI}}$ denotes the probability of a transition from an arbitrary major chord to another major chord, where the respective root notes have an interval of four semitones upwards. This way we obtain transposition invariance of our features.

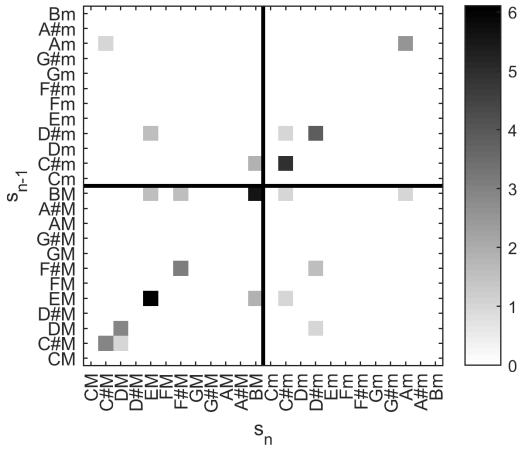
In the matrix A^{TPI}, the self-transition frequencies depend on the tempo of the recording. The self-transition probabilities might also contain valuable information about the recording but this is not connected to the harmonic content. Therefore we discard their effect by setting all elements on the main diagonals of the local transition matrices to zero and re-normalizing the matrices, such that the overall sum is one again. That way we assure that the tempo has no influence on the features. After the modifications, the matrix is now highly redundant. We need to discard several entries. First, we only need one column of each chord-type sub-matrix, because of the transposition invariance. Also we need to discard the self-transition probabilities, which are always zero and therefore contain no information. In the end, we use only a small subset of the matrix, which comprises all independent and relevant values as feature.

In Figure 3.12, we show the whole process of the feature generation using the Bach example. For this example, we use the Viterbi algorithm since the results are visually better understandable. At first, we see a chord estimation result with the respective chord transitions shown as red

(a) Estimated states and transitions



(b) Chord transition matrix A^{w}



(c) TPI chord transition matrix A^{TPI}

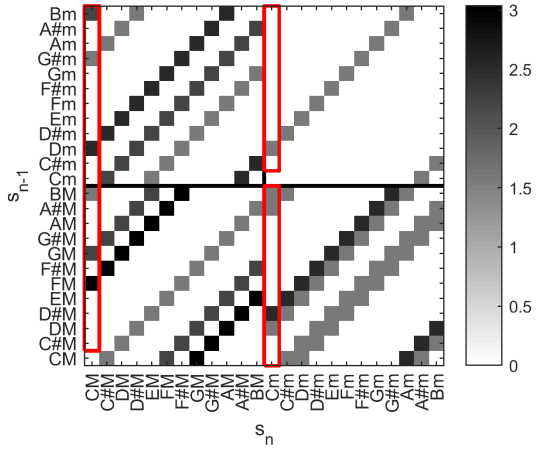


Figure 3.12: The Viterbi-based feature generating process performed on the Bach example. (a) shows the estimated state sequence and the corresponding state transition. (b) depicts the the chord transition matrix and (c) shows the TPI transition matrix without the main diagonal elements. The marked elements are statistically independent and serve as feature vector. We display all matrices after logarithmic compression ($\eta = 1000$, similar to Equation 2.10).

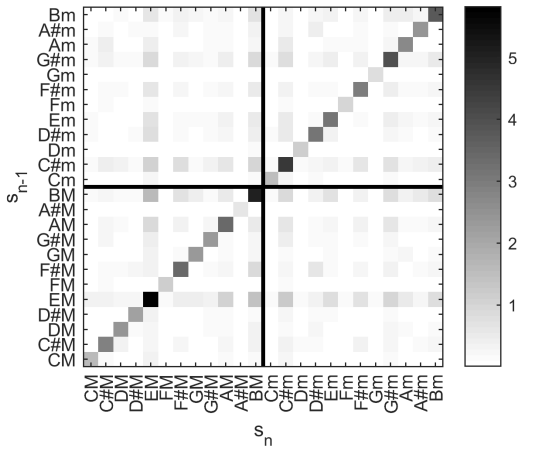
lines between the chords. By averaging the transitions over time, we arrive at the depicted chord transition matrix, see Figure 3.12b. We see that the matrix is sparse which shows that there are only few distinct types of transitions detected. The biggest values occur on the main diagonal, which contains the self-transition probabilities. The values on the main diagonal are a measure for the time the respective chords are played. We see that among the self-transitions, the chords EM, and BM have the highest self-transition probabilities, which is consistent with the recognized chords. For clarity, the plot is divided into four sub-matrices which separate transitions between pairs of chord types (black lines). For the next matrix, we perform the TPI transform and set the self-transition probabilities to zero. We see that the elements on each diagonal of the sub-matrices are now equal. Therefore, we have a high redundancy in this matrix. The entries marked with a red frame show the relevant information which is not redundant. Therefore, we select those entries for the final feature vector. That way, we obtain a feature vector that should have the desired properties of invariance against tempo and key.

3.4.2 Influence of Double-Softmax Rescaling

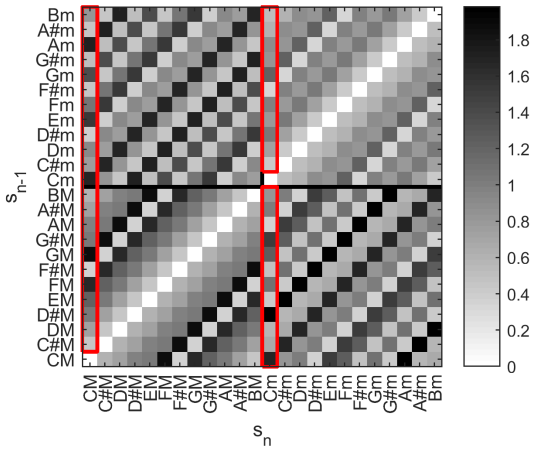
In Figure 3.13 a and b, we show the resulting matrices A^{w} and A^{TPI} for the Baum–Welch algorithm. Here, we directly applied the aggregation algorithm on $\xi_n(i,j)$. We see, that A^{w} is not sparse anymore as opposed to the Viterbi algorithm. For the most probable chord (EM) we observe that both transitions from and to this chord have a high probability. We see that the main diagonal still has the highest values. After the TPI transform we have strong diagonal structures. We see high values at rising and falling fifth transitions (e.g. CM $\rightarrow$ GM and CM $\rightarrow$ FM, respectively), which is similar to observations in [45]. However, we also see high values at transitions that we often find in the aforementioned lattice-structure (e.g. CM $\rightarrow$ Cm and CM $\rightarrow$ Em, compare Section 3.2.2), which are sometimes higher than transitions like the CM $\rightarrow$ Gm transition class, which we find in the estimated chord sequence and therefore assume to contain information about the harmonic progressions.

Previously we achieved good visual results by applying a double-softmax rescaling on $\xi_n(i,j)$. We were able to suppress noise-like transitions thoroughly. We now apply the aggregation algorithm on a rescaled $\xi_n(i,j)$, again with parameters $\eta_1 = 10$ and $\eta_2 = 10000$. We show the resulting matrices in Figure 3.13c and d. Since we delete the main diagonal in the rescaling process, it is not possible to show the self-transitions in A^{w} anymore. We see only a few dominant transitions over an almost uniform background. Comparing them with the result from the Viterbi algorithm in Figure 3.12, we observe a high similarity. However, we see some additional transitions, like EM $\rightarrow$ AM, which we already discussed previously. We see that we were able to obtain additional information without capturing the noise-like transitions. On the other hand, we apply a strongly non-linear operation, which is very sensitive on the input and which may yield very different

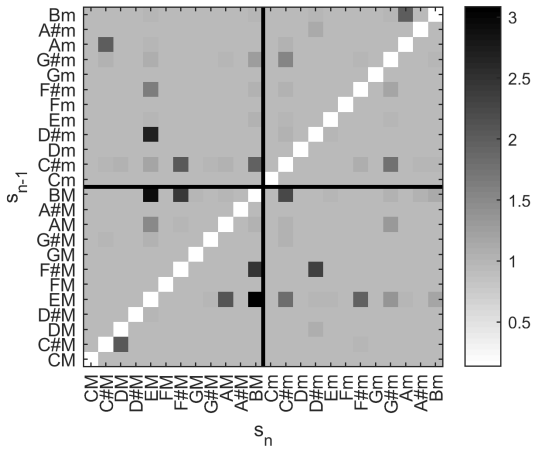
(a) Chord transition matrix A^{w}



(b) TPI chord transition matrix A^{TPI}



(c) A^{w} (double-softmax)



(d) A^{TPI} (double-softmax)

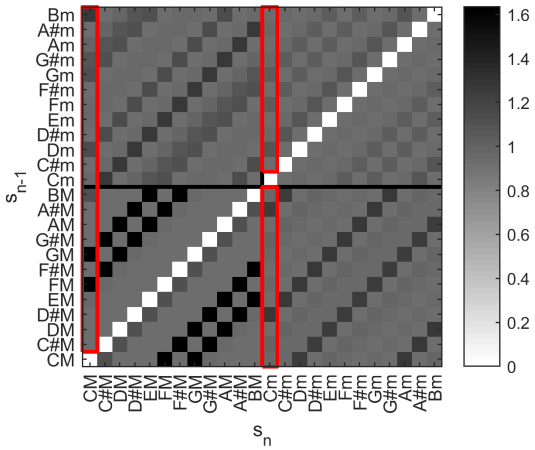


Figure 3.13: The chord transitions matrices of the Bach example, generated with the Baum–Welch-based approach. We show the original matrix in (a), the transposition invariant matrix in (b). Here, we marked the elements of the resulting feature vector. Sub-figures (c) and (d) show the same matrices when we use a double-softmax rescaling before the aggregation. We display all matrices after logarithmic compression ($\eta = 1000$, similar to Equation 2.10).

behavior on different pieces. It is therefore questionable how suitable the double-softmax-rescaled feature performs in a classification scenario.

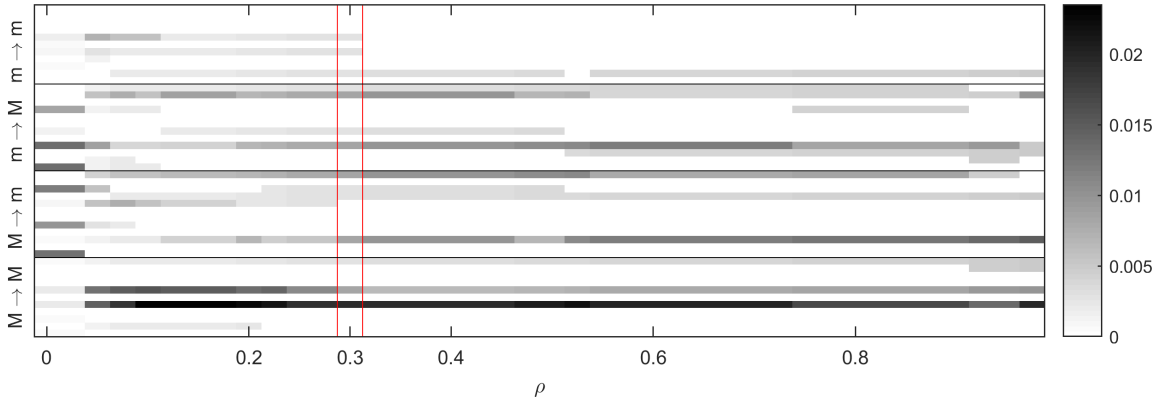
3.4.3 Influence of the Self-Transition Probability

In Section 3.3, we examine the influence of the self-transition probability ρ on the chord recognition result and the local chord transition estimation. Now, we take a closer look at the derived piece-level features for different values of ρ. For this experiment, we generate piece-level features of the Bach example with both approaches using different values of ρ. Figure 3.14 shows the resulting feature vectors. The horizontal axis shows the different values of ρ. The column above each value represents the corresponding feature vector as intensity plot. The vertical axis is split in regions that represent transitions between different chord modes. Within each region, the values are ordered in ascending order by the size of the interval of the respective transition. Since we excluded self-transitions, each transition type that does not change the chord type has only 11 entries, whereas the remaining transition types have 12 entries. If we examine the progression with increasing ρ, we see a smooth behavior for the Baum–Welch algorithm as opposed to sudden changes for the Viterbi algorithm. This is due to the discrete transition counting. There are only few transitions taken into account (the chord is expected to change around 30 times in this recording), and the change of one chord label changes two transitions. Thus, the change of a single chord has a big impact on the overall distribution of the different transitions.

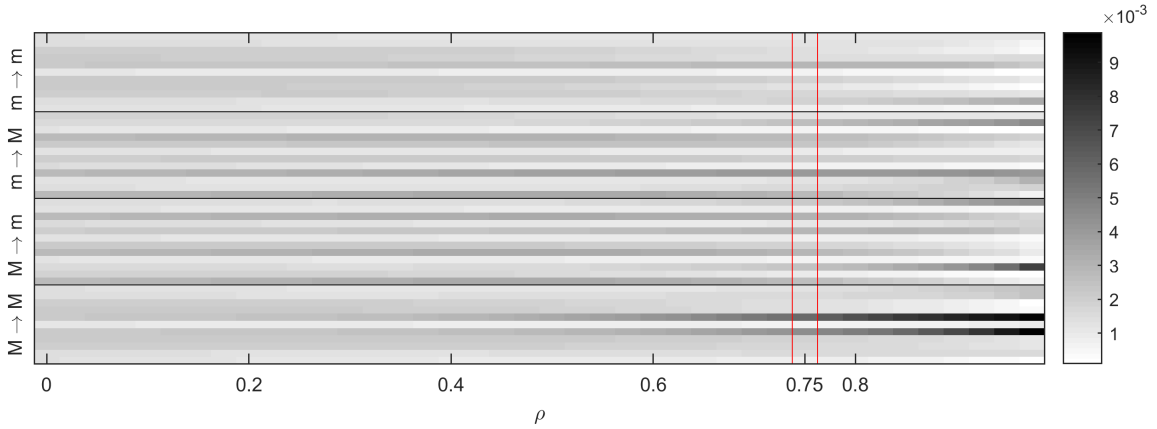
For $\rho < 0.05$, we see a different behavior. Here, we observe only a few transition types, which do not frequently occur for bigger values of ρ. Those transitions occur between chords that have two notes in common. A template-based chord recognition algorithm is likely to confuse those transitions. As explained in Section 3.3, for small values of ρ, the Viterbi-algorithm changes back and forth between the two most common chords. These transitions are much more frequent than the others, which is why they dominate here. This behavior stops abruptly as soon as the self-transition probability rises above the other non-self-transition probabilities. On the other side, we see that the feature vector gets sparse with increasing ρ, because there are fewer chord changes, thus containing less information. Since both extreme values of ρ are apparently not suitable features, there must be an optimal ρ in between. The vector obtained with standard $\rho = 0.3$ is marked red and lies in the supposedly good region, where we assume semantically meaningful feature vectors.

The Baum–Welch-based features (see Figure 3.14b) have a nearly uniform distribution for small values of ρ and become more similar to the Viterbi-based features for large values of ρ. The middle area ($0.4 < \rho < 0.8$)has a transition between those extremes. Again, one can argue that the feature vector gets sparser for large $\rho > 0.9$ thus carrying less information. On the other hand, the feature vector becomes more uniform with small $\rho < 0.2$, which also seems to carry

(a) Viterbi



(b) Baum–Welch



(c) Baum–Welch (double-softmax)

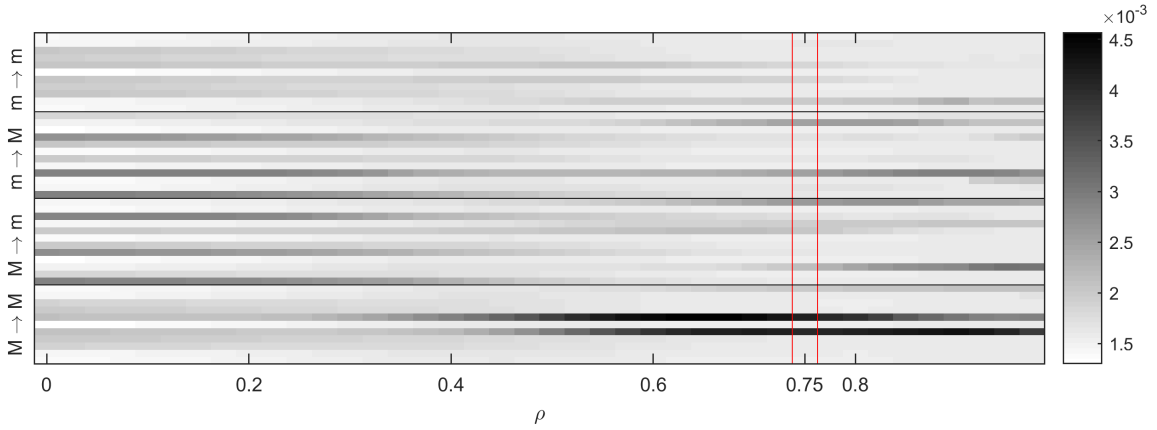


Figure 3.14: Chord transition features generated with the (a) Viterbi, (b) Baum–Welch algorithm and (c) Baum–Welch algorithm with double-softmax rescaling for different ρ. The optimized ρ is marked with a red frame.

less information. By these considerations, the optimal ρ must be somewhere in between. For the Bach choral, the value $\rho = 0.75$—as optimized on the *Beatles* dataset—seems to be a good compromise. We also see, that the vector is not very sensitive to small changes of ρ. Thus, small deviations should still yield similar results.

If we perform double-softmax rescaling before the feature aggregation, we obtain a different behavior (see Figure 3.14c). Around standard ρ, we clearly see the relevant transitions. When we decrease ρ, the vector becomes sparser and "lattice" transitions become dominant. We have to keep in mind that the softmax parameters were visually optimized on a chord transition estimation result with $\rho = 0.75$. Therefore, we cannot expect the algorithm to work for other values of ρ. Note that those arguments are purely qualitative and based on only one example. We perform a more systematic examination in Chapter 5.

Chapter 4

Background: Style Classification

In this chapter, we introduce the problem of musical style classification. Knowledge of a musical style can be important in tasks like music retrieval or to categorize a music library. To classify the style of a recording, we need a good classifier as well as suitable features. We covered our proposed feature generation algorithms in the previous chapter. In this chapter, we introduce the remaining components of the problem. At first, we present the musical scenario of our experiments. We describe the dataset and provide some background information to further specify the problem. Afterwards, we briefly cover basic machine learning techniques, needed for the classification. In the last section, we give a small overview over previous work in this area.

4.1 Musical Scenario

To systematically evaluate style classification systems, we need a good dataset that contains recordings of different styles. In [45], Weiß introduced the *CrossEra* dataset for this purpose. The dataset comprises four different sub-styles of Western classical music from 1650 until the present. The dataset splits into four classes: Baroque, Classical, Romantic, and Modern. Therefore the pieces cover four important historical periods or "eras". Each style is represented by 400 recordings. Of those 400 recordings, there are each 200 piano and 200 orchestra pieces. For some experiments, we need the subsets *CrossEra-Piano* and *CrossEra-Orchestra*, which consist of only piano and orchestra pieces, respectively. Table 4.1 shows the content of the *CrossEra* database, grouped by style and composer. We see that the database is balanced with respect to styles and instrumentation but not with respect to the composers. Unlike pop music, there are many different versions of one composition or work in classical music. Especially the popular works have been performed and recorded many times. Each recording differs from the others in several ways, including tempo, possibly key, specific instruments and recording equipment. The *CrossEra* database mainly contains recordings from the Naxos label. In *CrossEra-Piano*,

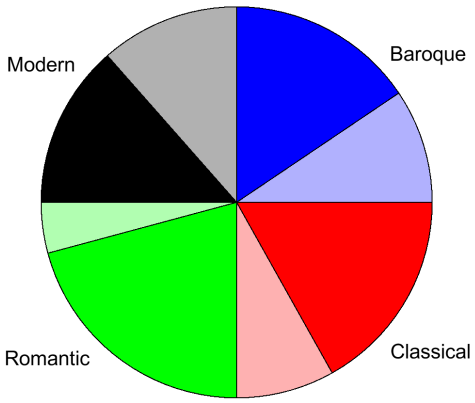


Figure 4.1: Number of recordings in the *CrossEraMirror* database in relation to the recordings in CrossEra

Weiß only selected recordings performed on a modern grand piano [45]. This excludes the harpsichord, which was commonly used in the Baroque era. This measure prevents a classification based on the instrumentation. We perform most of our experiments on the *CrossEra* set. In the end we validate our results using unseen data. For this purpose, we use another dataset, called *CrossEraMirror*. As the name indicates, it comprises different recordings of the same works as *CrossEra*. With this dataset we can perform cross-version experiments. Not every recording in *CrossEra* has an equivalent in *CrossEraMirror* interpreting the same work. The content of *CrossEraMirror* is also listed in Table 4.1. To visualize the difference between the two databases, Figure 4.1 shows a pie chart with the distribution of the different eras in the database. Each quadrant represents one era in the *CrossEra* database. The percentage which is also in *CrossEraMirror* is colored fully, the part which is only in *CrossEra* is colored transparently. We see that *CrossEraMirror* is no longer balanced with respect to era, since Modern and Baroque are underrepresented. The dataset is also not balanced with respect to instrumentation.

4.2 Basic Machine Learning Concepts

In this section, we give a brief introduction of concepts connected with machine learning and classification experiments. This is a wide field so we limit ourselves to topics relevant for this thesis. At first, we explain the basic notions of machine learning—also known as pattern recognition—

Style	Composer	Number of Recordings	
		CrossEra	*CrossEraMirror*
Baroque		400	249
	Albinoni, Tomaso	23	3
	Bach, Johann Sebastian	121	121
	Corelli, Arcangelo	15	14
	Couperin, Francois	36	36
	Giustini, Lodovico	8	0
	Handel, George Frederic	38	16
	Lully, Jean-Baptiste	15	1
	Platti, Giovanni Benedetto	41	0
	Purcell, Henry	10	4
	Rameau, Jean-Phillipe	60	41
	Vivaldi, Antonio	33	13
Classical		400	270
	Bach, Johann Christian	30	11
	Boccherini, Luigi Rodolfo	23	23
	Cimarosa, Domenico	32	1
	Clementi, Muzio	33	31
	Dussek, Jan Ladislav	35	0
	Haydn, Johann Michael	11	0
	Haydn, Joseph	100	95
	Mozart, Wolfgang Amadeus	114	109
	Pleyel, Ignace Joseph	10	0
	Salieri, Antonio	12	0
Romantic		400	333
	Berlioz, Hector	8	8
	Borodin, Alexander	9	9
	Brahms, Johannes	37	37
	Bruckner, Anton	12	12
	Chopin, Frederic	29	29
	Dvorak, Antonin	12	12
	Faure, Gabriel	15	13
	Grieg, Edvard	40	40
	Liszt, Franz	34	25
	Mendelssohn-Bartholdy, Felix	48	25
	Mussorgsky, Modest	12	0
	Rimsky-Korsakov, Nicolai	7	4
	Saint-Saens, Camille	16	8
	Schumann, Clara	13	0
	Schumann, Robert	43	43
	Smetana, Bedrich	6	6
	Tchaikovsky, Peter Iljitsch	30	19
	Verdi, Giuseppe	20	20
	Wagner, Richard	9	9
Modern		400	216
	Antheil, George	8	0
	Bartok, Bela	29	20
	Berg, Alban	11	10
	Boulez, Pierre	11	0
	Britten, Benjamin	9	9
	Hindemith, Paul	38	28
	Ives, Charles Edward	6	0
	Messiaen, Olivier	42	0
	Milhaud, Darius	22	0
	Prokofiew, Sergej	37	12
	Schoenberg, Arnold	30	18
	Shostakovich, Dmitri	82	82
	Stravinsky, Igor	28	23
	Varese, Edgar	14	0
	Webern, Anton	17	5
	Weill, Kurt	16	0

Table 4.1: Content of the *CrossEra* and *CrossEraMirror* database grouped by style and composer.

before we move on to the concept of cross validation, explain dimensionality reduction, and describe the Gaussian classifier.

4.2.1 Basic Notions

Machine learning systems generate models which are able to predict the outcome of experiments [4]. The key aspect is that these models are not generated by a human but by a machine. The machine is able to classify data using knowledge it gained from previous data. We use this concept for our classification experiments by training a model for style classification based on chord transition features. Each machine learning experiment consists of a *training phase* and a *test phase*. In the training phase we train the model using one part of the available data. In the test phase we classify the remaining data based on the trained model.

We distinguish supervised and unsupervised training methods. In *supervised training*, the system knows the ground truth labels of the training data. *Unsupervised training* only uses unlabeled data. A good example for unsupervised training is the Baum–Welch algorithm for HMM training, since the true internal states of the HMM are not needed. Learning the Gaussian emissions from labeled data as performed in Section 2.3 fall in the category of supervised methods [40].

Our system consists of three steps. The feature extraction stage transforms the raw data into a new format which is suitable for the classification task. This part is covered in Chapter 3. Dimensionality reduction reduces the number of elements of the features vector by finding appropriate combinations of the feature components that have a high discriminative power. This way we can discard irrelevant information. Finally, we use a classifier to compute the estimated labels from the reduced features. The dimensionality reduction and the classifier require a training phase.

For the classification to work, the features need to have a property called *separability*. The features belonging to one class should be close to each other (low intra-class distance) and the features of different classes should be far from each other (high inter-class distance) [15].

4.2.2 Cross Validation

Any experiment on machine learning systems consists of a training phase and a test phase. In the training phase, we use data to train a model until it yields an optimal outcome. Now, the system performs well on the training data, but there is not guarantee, that it performs well on previously unseen data. The model might be too specific to generalize well. This problem is known as overfitting a system and describes the loss of generality of the learned patterns [7]. Therefore, it is very important to evaluate the system only on data that was not used for training. This way we are close to a real world scenario where we do not know the labels of the test data.

Figure 4.2: Example of how a database can be split three times into test an training set in order to perform a 3-fold cross validation.

It is common practice to split the database in two sets, the training set and the test set. Each set is only used for training and testing, respectively. The results reported after the experiment come only from the test set. To still use the entire dataset for evaluation a so-called *cross validation* [38] is used. We split the whole dataset in n so-called *folds* and perform n experiments. In each experiment we use one fold as test set and the remaining folds for training. In this case, we talk of an n-fold cross validation. That way we can evaluate the classification system on all elements of the set. Figure 4.2 shows an example of how a database can be split for a 3-fold cross validation.

There are different constraints on how to perform the split. A very common one is *stratification* [28]. We call a set stratified if it contains an equal number of samples of each class. It makes sense to stratify both the test and the training set. This can be achieved by successively assigning randomly chosen samples from one class to each fold. Sometimes the structure of the dataset does not allow perfect stratification. In this case we can either randomly delete or duplicate samples within the set until the set is stratified.

In some datasets, there exist groups of samples that belong to the same class but have other similarities. For example two recordings from the same album[1] were probably recorded with the same equipment and by the same people, with the same instruments. The characteristics of each instrument, musician and equipment can contribute to the features. A certain recording engineer for example could tend to suppress the high frequencies slightly stronger than their colleagues, which has an impact on the features. A classifier might now use that information to determine that two recordings come from the same album, thus inferring that they have the same style. While this decision is probably correct, the classifier bases this decision on properties which are not related to the musical style. The classifier would not work as well on other recordings of the

[1]In the context of classical music, the term "album" refers to a group of recordings which were recorded together, i.e. in the same recording session or live recordings of the same concert.

same style processed by a different recording engineer. In the MIR society, this phenomenon is known as the "*album effect*" [12, 47]. The album effect is one type of overfitting.

To avoid the album effect, we have to make sure that the whole album is either in the training or the test set. That way we increase the probability of overfitting to an album. But, since we excluded all samples which profit from the overitted classifier from the test set, the classifier cannot use recording artifacts for the classification. We call this procedure an "*album filter*". Using such filters typically makes the results worse. But in this case, we benefit from a better approximation to a real world scenario. Since the album annotations are incomplete in our dataset, we instead introduce an even stronger restriction: the *composer filter*. Here, we make sure that works from each composer only appear in either the training or the test set. The composer filter includes an album filter, since each album in the database only contains compositions from one composer. By this measure we also avoid influences of overfitting on a certain composer style.

4.2.3 Dimensionality Reduction

Dimensionality reduction is used to further compress the features. A high dimensional feature vector has several disadvantages, which are summarized as the *curse of dimensionality* [40, 42]. In general, we can say, that a system with high dimensional features is more prone to overfitting and generalizes worse. However, low-dimensional features may carry less information. Thus the parameters and strategy of the dimensionality reduction play a crucial role for the overall result.

We explain two basic algorithms: the principal component analysis (PCA) and linear discriminant analysis (LDA). PCA is an unsupervised algorithm while LDA is supervised but, therefore, also prone to overfitting.

4.2.3.1 Principal Component Analysis

This chapter is a brief summary of [40, Chapter 6.3], which gives a much more detailed introduction. *Principal component analysis*, also known as Karhunen–Loève transform serves two purposes. First, the transform performs a linear operation on the feature vectors, such that all elements become uncorrelated. The basis vectors of the new feature space are the principal components of the training set, giving this method its name. The transform does not change the dimensionality. However, now that the elements of the feature vector are decorrelated, we can easily compute the variance of each element and discard the elements with the lowest variance, since they probably have the smallest influence on the separability of the classes.

The principal components are the eigenvectors of the correlation matrix $R_{\mathbf{f}}$, which can be estimated on a training set of P feature vectors $\mathbf{f}_p$, $p \in [1:P]$:

$$R_{\mathbf{f}} = \frac{1}{P} \sum_{p=1}^{P} \mathbf{f}_p \mathbf{f}_p^{\intercal}. \tag{4.1}$$

We can now compute the eigenvectors $\mathbf{a}_i$ in descending order with respect to the corresponding eigenvalues λ_i. We summarize the eigenvectors in the transform matrix A_{PCA}. We can write the transformed feature vector $\tilde{\mathbf{f}}$ as:

$$\tilde{\mathbf{f}} = A_{\text{PCA}}^{\intercal} \mathbf{f} \tag{4.2}$$

with the correlation matrix

$$R_{\tilde{\mathbf{f}}} = A_{\text{PCA}}^{\intercal} R_{\mathbf{f}} A_{\text{PCA}} = \Lambda. \tag{4.3}$$

Here Λ denotes the eigenvalue matrix, a diagonal matrix with the eigenvalues as elements in descending order along the main diagonal. We see from this equality that the eigenvalues of $R_{\mathbf{f}}$ are equal to the variances of the feature elements in the new vector space. To reduce the dimensionality of the feature space from P to $Q \leq P$, we can take the first Q elements of the transformed feature vector. We therefore define $A_{\text{PCA}} \in \mathbb{R}^{P \times Q}$ as the matrix containing the eigenvectors corresponding to the k highest eigenvalues.

However, since the PCA does not take the class labels into account, the dimensionality reduction might delete important information. It is possible that two classes are only separable in the dimensions with low variance. In Figure 4.3 we show an example of PCA leading to a loss of separability. We see that the two classes are linearly separable in the two dimensional plane. Performing a PCA we obtain two principal components $\tilde{f}_1$ and $\tilde{f}_2$. When we project the set on $\tilde{f}_1$ both classes overlap. The set is not separable anymore. This example shows, that we must use PCA with caution in order to not destroy important information.

4.2.3.2 Linear Discriminant Analysis

In contrast to principal component analysis, linear discriminant analysis is a supervised procedure. Here, we jointly maximize the inter-class variance and minimize the intra-class variance. In the new feature space the classes are compact and have a high distance from each other. That way we keep maximal separability.

Again, we search for a matrix A_{LDA} such that

$$\tilde{\mathbf{f}} = A_{\text{LDA}}^{\intercal} \mathbf{f} \tag{4.4}$$

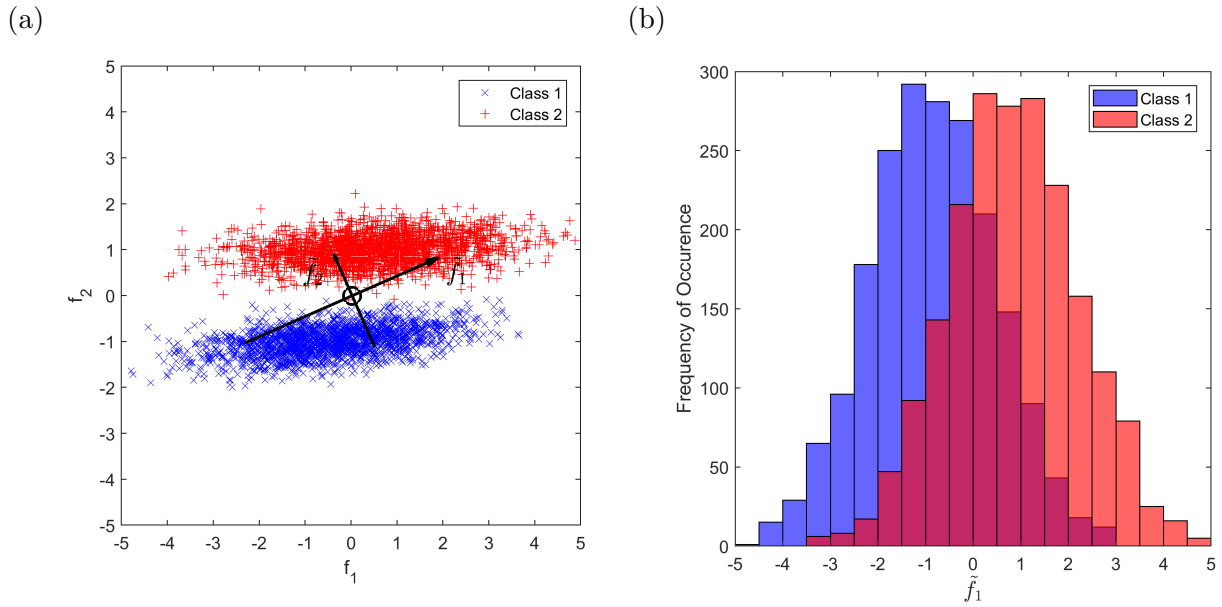


Figure 4.3: This figure shows a dataset in two dimensions, which is processed with PCA (a). The principal components are drawn as a second set of axes. The length of the axes indicates the variance of the dataset along this axis. Sub-figure (b) shows the distribution of the data points with respect to $\tilde{f}_1$.

has the desired properties. Using Fisher's criterion [11] as described in [40], the basis of this vector space are the eigenvalues of $S_w^{-1}S_b$, with the intra-class scatter matrix $S_w \in \mathbb{R}^{P\times P}$ and the inter-class scatter matrix $S_b \in \mathbb{R}^{P\times P}$, with P as the size of the feature vectors. It can be shown that for a C-class problem it is sufficient to choose the vectors corresponding to the $C-1$ highest eigenvalues, since all other components would be linearly dependent to the first $C-1$ components. In [40, Section 5.8], Theodoridis explains LDA in more detail.

Figure 4.4 shows the same dataset as in the previous section. This time we process the data with LDA. The set is still separable in the two dimensional space. We draw the projection axes LDA computed. We see that after the projection, most elements of class 1 are on the negative range of the $\tilde{\mathbf{f}}$ axis whereas most class 2 elements are on the positive range. The classes are still separable after we projected the data on $\tilde{f}_1$. The histogram for the reduced data confirms this, as we see that the classes do not overlap.

In [40] Theodoridis states that the LDA does not perform well on very high-dimensional data. Particularly if the dimensionality exceeds the number of training samples per class, S_w is not invertible anymore. Theodoridis suggests to combine PCA and LDA. In this approach, we first reduce the dimensionality with PCA. Afterwards, we train LDA on the reduced vectors and

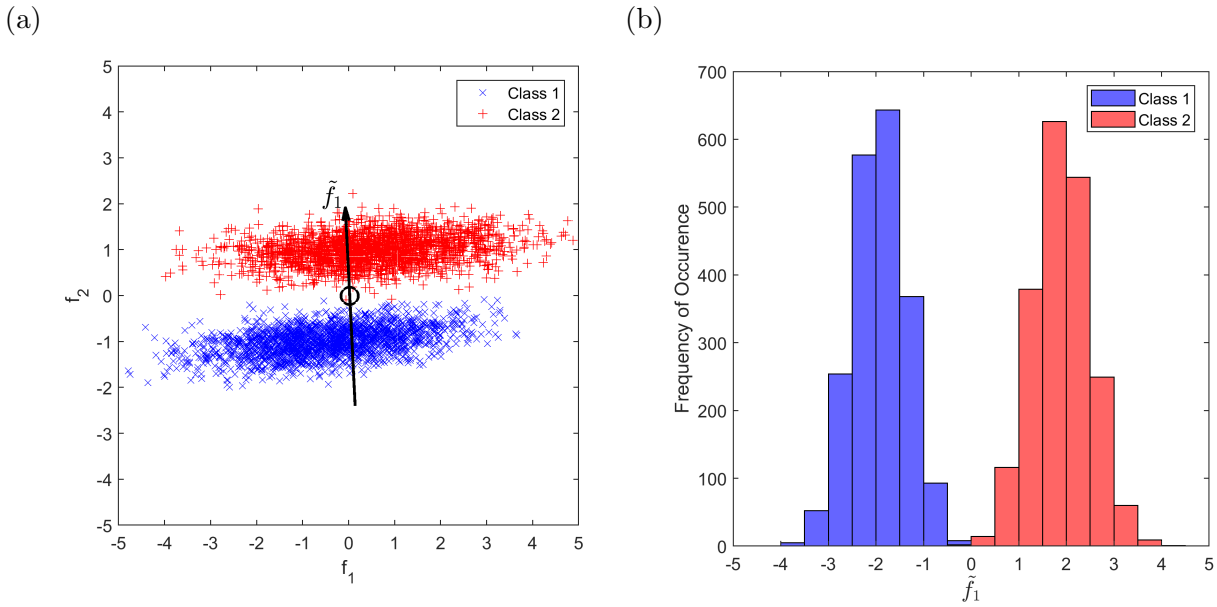


Figure 4.4: This figure shows a 2D-dataset which is processed with LDA (a). The basis vector on the reduced feature space is drawn as an arrow. Sub-figure (b) shows the distribution in this space.

perform LDA to further reduce the dimensionality. We define the reduced feature vector $\tilde{\mathbf{f}}$ as follows:

$$\tilde{\mathbf{f}} = A_{\text{LDA}}^{\intercal} A_{\text{PCA}}^{\intercal} \mathbf{f}. \tag{4.5}$$

4.2.4 Gaussian Classifier

After we obtained dimensionality-reduced feature vectors, we need to use a classifier to estimate the class label. In this thesis we use a Bayesian classifier based on a Gaussian mixture model (compare [40, Chapter 2]). A Gaussian mixture model is a superposition of several Gaussian PDFs. We train a model for each class that describes the distribution of the respective data points. For each class we use the training data of that respective class to learn the model parameters with the expectation-maximization algorithm (compare [40, Section 2.5.5]). We use the training method implemented in the *scikit-learn*[2] [28] toolbox. Afterwards, we evaluate the models from all classes for each instance from the test set. We choose the class, whose model yielded the highest probability as class label. In this thesis, we only use a single Gaussian PDF, since the results from [45] suggest that one Gaussian PDF is sufficient. Using several Gaussian PDFs in the mixture model increases the number of model parameters. Therefore, the system is more prone to overfitting.

[2] `http://scikit-learn.org`

4.3 Related Work

This section provides an overview over related work in style classification. A musical style can roughly defined by certain characteristics regarding instrumentation, melodic and harmonic content, dynamics and timbre [18]. Assigning a piece to a style category can be ambiguous since styles are not always well-defined, can possible overlap, and can be further divided [13]. Therefore problem of style classification is an equally widespread problem consisting of many sub-problems. We can distinguish between different genres, such as Jazz, Popular and Classical [29,30] or Chinese folk music and Western popular music [35]. Some genres can be split in different sub-styles, which poses the problem of sub-style classification. Here, we distinguish between sub-styles belonging to the same genre. One example is our problem of era classification within Western classical music [45]. Another example is the recognition of Jazz sub-styles in [19]. We can go one step further and classify different composer within a (sub-)style. In [45], Weiß performs experiments to distinguish between several Western classical composers. In [19], Hedges *et al.* classify different Jazz composers.

Not only the set of styles varies within the problem of style classification, also the available information for the feature calculation may be different. We can distinguish between symbolic features and audio features. Symbolic features are based on symbolic music representations, like sheet music or MIDI (Musical Instrument Digital Interface) files. In such approaches we can directly use the musical information. In [31], Ponce de León and Iñesta propose a set of MIDI-based features that combine melodic, harmonic and rhythmic features. Armentano *et al.* use MIDI-based features which describe sequences of MIDI events [1]. Ogihara and Li propose MIDI-based chord progression features [27], by combining several chords to n-grams.

On the other hand, audio features are always based on a recording. One approach is the use of standard audio features, like for example spectral envelope, zero crossing, or spectral flatness measure. A common choice is also the mel frequency cepstral coefficient, which is also useful in the field of speech recognition [41]. In [45], Weiß evaluates many of these features on the problem of sub-style classification of Western classical music. While the features perform well without any cross validation filter, the performance drops considerably when Weiß applies the composer filter. Another possibility is the use of *tonal features*. Tonal features are designed to contain musical information, such as features describing melody [34] or the frequency of chord transitions [45]. We call this kind of features, which have an explicit meaning, *high-level features* [45]. In [30], Pérez-Sancho *et al.* interpret the feature generation process as transforming the audio data to a symbolic representation. On the other hand, there are *mid-level features*. In [45], Weiß defines mid-level features as feature that "relate to human-interpretable concepts but in a way that is not obvious." He shows that tonal mid-level perform well in the context of sub-style classification of Wester classical music.

In [45], Weiß proposes features that describe the occurrence of certain intervals and chord types. Weiß uses chroma templates describing each interval and each triad. The feature type is based on the similarities between each chroma vector and each template vector. The feature vector comprises the average and the standard deviation of those similarity measures across the recording. Weiß also proposes features which quantify tonal complexity, referring to the distribution of pitch class energies [47]. The feature consist of various complexity measures quantifying for example sparsity of the chroma vector (low complexity) or the consonance (low complexity) of corresponding notes.

Additionally Weiß introduces high-level chord transition features [45]. The feature requires estimated chord labels. Weiß uses the *Chordino* system to estimate the chords with a large chord vocabulary. However, the feature aggregation algorithm only considers transitions between major and minor chords. In the end, the chord transition features contain information about the relative frequencies of the chord transitions.

In [45] Weiß showed, that tonal features work well in a style classification scenario. The template-based features showed the best performance. However, also the chord transition features yielded good results. Both feature types performed well also with composer filter. We can interpret our mid-level chord transition features as a conceptual combination of high-level chord transition features and the mid-level template based features. The features proposed in [45] constitute the baseline for the experiments in the following chapter. For a fair comparison we apply the same classification pipeline on these and on our features.

Chapter 5

Style Classification Experiments

In this thesis, we focus on sub-style classification within Western classical music. In particular we work on the *CrossEra* dataset described in Section 4.1. In this chapter, we present the style classification experiments. At first, we introduce the classification pipeline, using the elements we describe in Section 4.2. We start with baseline experiments by evaluating some of the features described in [45] using our classification pipeline. Next, we give a comparison of the Viterbi-based and the Baum–Welch-based features and test combinations of chord transition features, template features, and complexity features. In the next experiments, we examine the influence of the self-transition probability in the HMM on the classification results and also the combination of different features. Furthermore, we look at the rescaling of the local chord transition features (compare Section 3.2.2), which improve the visual results greatly. We then demonstrate the album effect and examine how well different features generalize over different composers and instrumentation. The chapter is concluded by cross-version experiments to verify the previous results using the *CrossEraMirror* database. In this chapter, we give a detailed description of the experiments. In Appendix B.3, we give summarized results, which extend the results presented in this chapter.

5.1 Classification Pipeline

In the following experiments, we consistently use a specific processing pipeline to obtain our classification results. We summarize the pipeline in Figure 5.1. We give an explanation of all steps in the following. First, we divide the data set into a training set and a test set. We perform the split employing a composer filter as described in Section 4.2.2 and forcing stratification by randomly removing elements. We use about two thirds of the dataset for training and one third for testing. In the next step, we extract piece-level features from the recordings of both

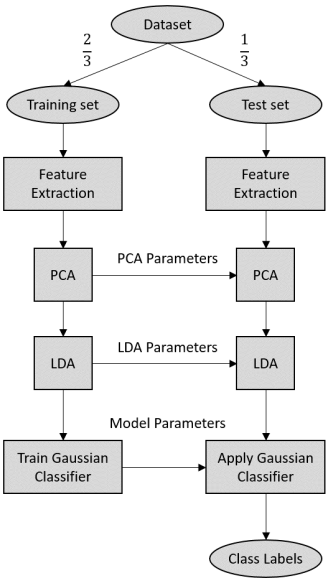


Figure 5.1: Visualization of the classification pipeline.

sets. We have already introduced the algorithms in Chapter 3. For comparison we also use the template-based features from [46] and the complexity features from [47].

For dimensionality reduction we use PCA (Section 4.2.3.1) followed by LDA (Section 4.2.3.2). First, we learn the PCA parameters on the training set. We then perform PCA using those parameters on both sets. We reduce the feature dimensions to $Q = 200$ if the dimensionality is greater than 200 and keep it constant if the feature vector has less than or exactly 200 dimensions. If we perform experiments on the subsets *CrossEra-Piano* or *CrossEra-Orchestra*, we set this threshold to 100 to account for the smaller training set. We need to use a dimensionality reduction before LDA to avoid overfitting LDA. We found that the result is not sensitive on the exact choice of this parameter as long as we choose Q smaller than the number of training samples but not too small such that we delete important information. Setting the parameter Q to half the number of samples per class seems to be a reasonable choice. For small feature vectors, we do not need the PCA but for reasons of consistency, we always use it. In this case, we only decorrelate the elements of the feature vector. LDA is performed similarly. First, we learn the parameters on the training set and afterwards perform LDA on both sets. Since we are dealing with a 4-class problem, the resulting reduced feature vector has only three dimensions. As mentioned in Section 4.2.3.2, LDA can only produce $C - 1$ independent components for a C-class problem.

We use the reduced feature vectors of the training set to train a Gaussian classifier, which then classifies the test set. We compare the resulting estimated class labels to the annotation. We use the accuracy as quality measure, which is the ratio of correctly classified recordings. We perform a 3-fold cross validation (compare Section 4.2.2) such that we classify (almost) every recording

once. We repeat the cross validation ten times (ten "runs") and average the accuracy of the cross-validations to obtain a result that depends less on the specific splits. We make sure that the splits are exactly the same for every experiment we perform by setting a fixed seed for the random generators in each experiment. We also repeatedly tested random splits which yielded similar results as the fixed splits we use in this thesis. The resulting mean accuracy of the 10 runs is the main quality measure of the used setting. However, we consider two more measures, the inter-fold deviation and the inter-run deviation [39, 45]. The inter-run deviation indicates the stability over different splits. We average the accuracy over all three folds in each run. That way we obtain mean accuracies for all ten runs. Afterwards, we compute the standard deviation of these accuracies. On the other hand, the inter-fold deviation indicates the stability over different test and training data. We first compute the standard deviation of the accuracies over all three folds in each run. We obtain ten different standard deviations, which we then average. We use the public toolbox *scikit-learn* to implement the machine learning algorithms.

5.2 Baseline Experiments

The goal of this section is to establish a baseline by comparing our features to the features used for classification in [45]. We do not compare our results to the published results in detail. Instead we use the same features as input for our own classifier. That way, we can be sure that the circumstances are the same, since we use a slightly different classifier.

In this experiment we compare the features used in [45] to the proposed chord transition features using the a chord vocabulary consisting of the 24 different major and minor triads. We select the template-based features (denoted as "Template"), a combination of template-based and complexity features (denoted as "Complexity & Template") and the high-level chord transition features as baseline features. To compute our features we need to model an HMM (compare Section 2.4). We model the state transition matrix A with the method we discuss in Section 3.3. We use the standard values for ρ—we set $\rho = 0.3$ for Viterbi-based features and $\rho = 0.75$ for Baum-Welch-based features—as self-transition probability. We perform the experiments on the full dataset *CrossEra-Full*, as well as on the instrumentation subsets, *CrossEra-Piano* and *CrossEra-Orchestra*.

We give the results in Table 5.1. In this table we can see several interesting results. We see that the Viterbi-based features and the chord transition features from [45] perform similarly. This is according to our expectations, since we designed the Viterbi-based features with similar concepts. We see that the Template features from [45] and the Complexity & Template features perform better than all chord transition-based features. The distance is particularly high on the *CrossEra-Piano* set, where the difference is around 10%. This behavior is consistent with the results obtained in [45, Table 8.5].

Features	*CrossEra-Full*	*CrossEra-Piano*	*CrossEra-Orchestra*
Template	73.9%	74.0%	79.4%
Complexity & Template	74.7%	73.5%	79.3%
Chord Transitions (from [45])	67.2%	58.0%	72.1%
Chord Transitions (Viterbi)	68.7%	60.8%	74.3%
Chord Transitions (Baum–Welch)	69.3%	63.6%	75.8%

Table 5.1: Results of the baseline experiments. Different features from [45] and this thesis are compared. We perform the experiments on the full dataset *CrossEra-Full* as well as the subsets *CrossEra-Piano* and *CrossEra-Orchestra*.

Taking a look at the Baum–Welch-based features we see that they perform better than all other chord transition features. However, the distance to the Viterbi-based features is only 0.6% for the full *CrossEra* dataset. For the subsets the distance is bigger with 2.8% for *CrossEra-Piano* and 1.5% for *CrossEra-Orchestra*. We also see, that the performance on piano pieces is worse than on orchestra pieces. The piano pieces contain only a single instrument and therefore, have less variation in the timbre than the orchestra which contains many instruments. We find this behavior throughout all features, which is consistent with the results in [45]. We perform more experiments on the subsets in Section 5.8.

5.3 Influence of Different Chord Vocabularies

As a next step, we compare the results obtained with different chord vocabularies. We tested the *MajMin*, *Triads* and *Sevenths* chord vocabularies. The *MajMin* vocabulary contains all 24 major and minor triads. We extend this vocabulary with 4 augmented and 12 diminished triads to the *Triads* vocabulary, which contains 40 chords. The *Sevenths* vocabulary contains the 75 different seventh chords we introduce in Section 2.1.2. In Table 5.2 we give the obtained results. As first quality indicator, we use the obtained classification accuracy. Additionally, we give the inter-run deviation, which we use to get an idea of the magnitude of the difference.

We see that the Baum–Welch-based features and the Viterbi-based features perform similarly for the *MajMin* vocabulary. We observe different behaviors for features generated with the Viterbi and the Baum–Welch algorithm as soon as we change the chord vocabulary. The Viterbi-based features do not improve with increasing size of the chord vocabulary while the Baum–Welch-based features yield better results. Since we expect roughly 300 chord changes in an average recording of 10 minutes length, it is not possible to fully exploit all dimensions of high-dimensional feature vectors like the 464-dimensional feature vector of the *Sevenths* vocabulary. Instead many elements of the feature vector remain zero. We find the actual information only in a small part of the feature vector. The Baum–Welch algorithm on the other hand does not have this problem. Here, the underlying chord recognition scheme does not decide on one chord per time frame. Instead

		Viterbi		Baum–Welch	
Vocabulary	Dim.	Accuracy	Inter-Run Dev.	Accuracy	Inter-Run Dev.
MajMin	46	68.4%	0.6%	69.1%	0.8%
Triads	132	68.8%	0.4%	72.6%	0.6%
Seventh	464	67.8%	0.5%	74.6%	1.1%
Intervals	360	63.9%	0.9%	73.9%	0.7%

Table 5.2: Style classification results with different chord vocabularies. "Dim" denotes the original dimensionality of the feature vector as obtained from the feature aggregation algorithm.

we have the probabilities of all possible transitions. We are able to exploit ambiguities in the chord recognition process and are able to fully use the dimensionality of the feature vector. This leads to clearly improved results. The observed gain of 5.3 % between the *MajMin* and the *Sevenths* vocabulary is larger than the inter-run deviations.

Additionally, we define the *Intervals* vocabulary. This vocabulary comprises all 66 harmonic intervals that are distinguishable in the equal tempered scale respecting octave invariance and enharmonic equivalence. Note that we are not able to distinguish for example a major third over C and a minor sixth over E. Also, due to symmetry, we can only distinguish six tritone intervals. For example, the tritones over C and $F^\sharp$ are indistinguishable. This vocabulary has a special role in our system since it does not describe the chord transitions directly. However, every chord is comprised of several intervals. We can therefore interpret a transition between two chords as a superposition of several interval transitions. This is true for arbitrary chords such that we are able to implicitly describe a larger chord vocabulary. We now have a generalized feature which models all possible chords, avoiding problems like different numbers of notes per chord and HMMs with many states. Note that the HMM with the *Intervals* vocabulary has 66 states which is between the size of the *Triads* and *Sevenths* vocabularies. In Section 3.3, we observe a similar effect when we interpret a seventh chord as superposition of a major and a minor triad. We also see that this only works for the Baum–Welch algorithm, since the Viterbi algorithm decides on one of the possibilities. The same is true in this case. We therefore expect good performance for the Baum–Welch-based features but worse performance for the Viterbi-based features. The results shown in Table 5.2 confirm this hypothesis. The Baum–Welch-based features perform better than the *Triads* features and similarly to the *Sevenths* features, but the Viterbi-based features perform about 5% worse than any other chord vocabulary. The *Intervals* vocabulary shows the highest difference between Viterbi-based and Baum–Welch-based features.

We now present a more detailed analysis of the classification results. We start by looking at the confusion matrix of the classification experiment. Using the confusion matrix we can analyze the behavior of a classifier on a class level. Each entry of the matrix describes how often elements of one class are classified as the other class. The elements on the main diagonal are the so-called

per-class accuracies. Figure 5.2 shows the confusion matrix for three different features. The first matrix shows the results of the Baum–Welch-based features on the *Triads* vocabulary. There are several interesting points. We see that the class Modern has the highest accuracy. This is in line with musicology, since the rise of 12-tone music in this era has a very different chord structure. Among other differences, this distinguishes the Modern era clearly from the other eras. There are only very few confusions to the classes Baroque and Classical. About 10% of the recordings are misclassified as Romantic. This is an expected behavior since the two respective time periods are "neighbors" and have an overlap. In the earlier eras, we see many confusions between the classes Baroque, Classical and Romantic. Here, a confusion to the direct neighbor is not preferred, in fact confusions between Romantic and Baroque are more frequent than between Romantic and Classical. This is consistent with the results found in [45, Figure 7.6], which stated that some aspects of the chord transitions are more similar between the classes Baroque and Romantic than between Baroque and Classical. The class Romantic has the smallest accuracy, which constitutes a different behavior than observed in [45] with complexity and template-based features.

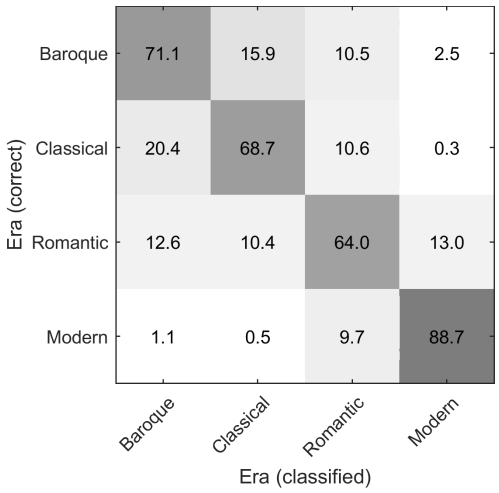
For the second matrix, we use the combination of complexity and template features from [45]. Here, we see the expected behavior, confusions to the "historically closer" eras are more common. However, there are now more confusions between Baroque and Classical.

5.4 Combination of Features Types

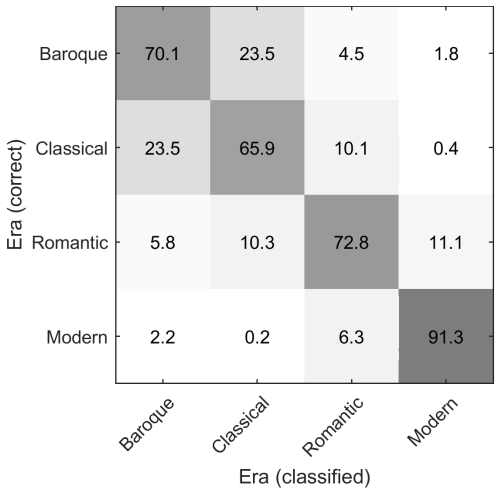
Since our Baum–Welch-based features are not able to clearly separate Romantic from Baroque but perform better in separating Baroque from Classical than complexity and template features, we now combine the two features. We show the result in the third matrix of Figure 5.2. We see that the per-class accuracies of Baroque, Classical, and Modern are larger than for either of the previous features. The accuracy of the class Romantic is larger than for the Baum–Welch-based features but slightly smaller than for the complexity and template features. We obtain an overall accuracy of 76.1% (compare Table 5.3), which is higher than the overall accuracies of each individual feature type alone. This shows that the information contained in both feature types is of different nature, such that each feature type benefits from the other.

We take this as a motivation to examine other combinations of feature types. Table 5.3 shows the classification accuracies of different feature combinations. We combine chord transition features with either template features, complexity features or a combination of both. As a reference, we also give the accuracies of the individual features alone (denoted with "None" in the respective other feature class). We see that complexity features alone perform worse than the template features, which agrees with the findings in [45]. We also see that a combination with Viterbi-based features is not always beneficial, similar to the findings in [45], where a combination of chord transition features with chroma-based complexity and template-based features is tested as well.

(a) Baum–Welch-based



(b) Complexity & Template



(c) Complexity & Template & Baum–Welch-based

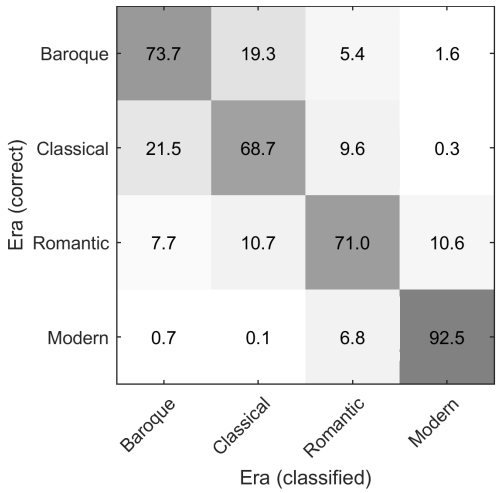


Figure 5.2: Confusion matrix of classification experiments with (a) Baum–Welch-based features, (b) complexity and template features (see [45]) and (c) a combination of both features.

Chroma-Based Features	Chord Transition Features		
	None	Viterbi	Baum–Welch
None	-	68.8%	72.6%
Template	73.9%	74.4%	76.0%
Complexity	64.4%	73.4%	75.9%
Complexity & Template	74.7%	74.0%	76.1%

Table 5.3: Classification accuracies of combinations of chord transition features and chroma-based features from [45].

We can also repeat the findings, that a combination of Viterbi-based features and template-based features performs best, better than the combination of Viterbi-based, Complexity, and Template. However, in [45] this difference was higher. The behavior of the Baum–Welch-based features differs from this. All combinations are better than each individual feature type. We obtain the best overall performance for these combinations. Interestingly all combinations also have very similar result. The differences of 0.1% are deemed insignificant. Apparently the amount of additional style information is the same for complexity and template features. This is remarkable, since complexity features alone have worse performance than template-based features. One possible interpretation is that complexity features are of different nature than the chord transition features. Therefore, there may be fewer overlap of the contained information. On the other hand, template-based features contain information about occurrence frequencies of different chord types. This information is related to our chord transition features. We can argue that much of the new information from the Baum–Welch-base features is redundant with the information contained in the template features.

In this section, we showed that—different to Viteri-based features—Baum–Welch-based contribute information to the tonal features from [45] and are able to further improve the classification accuracy up to 76%.

5.5 Influence of the Self-Transition Probability

In Section 3.4 we show the feature vectors of a single recording as a function of the self-transition probability ρ. We concluded that the standard ρ has a value where we assume a good quality. Those conclusions are only of qualitative nature since we did not yet have the possibility to test the features in a classification experiment. We now use the features derived with the *Triads* chord vocabulary using different values of ρ to classify the style of the recordings.[1] The experiment serves two purposes. First, we want to test if, a good chord recognition performance also leads

[1] We have to keep in mind that we obtain standard ρ by optimizing on a dataset consisting of modern pop music, which is different to Western classical music. Also the result depends on the annotations as described in Section 3.3.

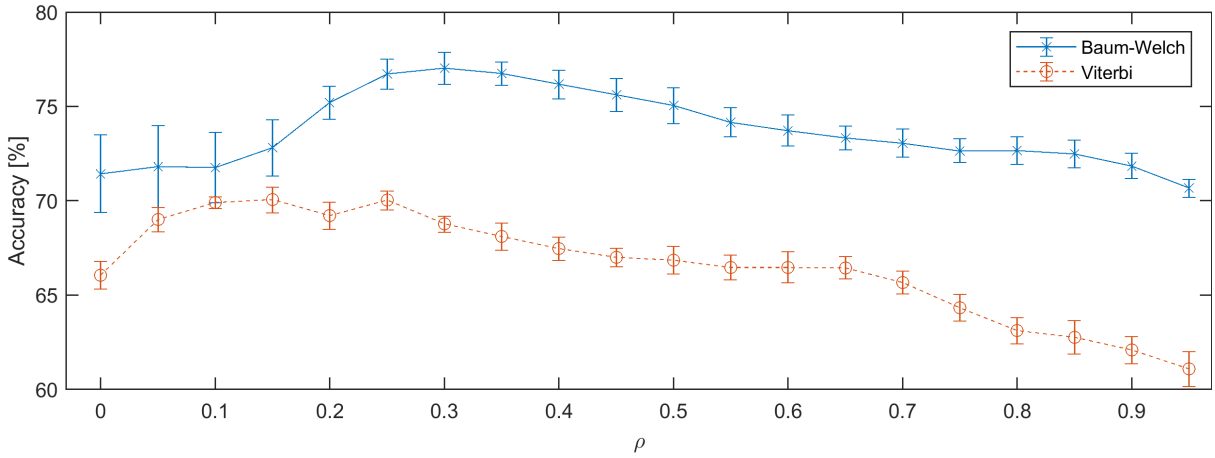


Figure 5.3: Classification accuracies for features generated with different ρ. The error bars indicate the inter-run deviation.

to efficient features. Second, we want to show that the gain of the Baum–Welch-based features compared to the Viterbi-based features is actually due to the algorithm and not due to the different choices of ρ.

In Figure 5.3, we show the classification accuracies of the Viterbi-based and the Baum–Welch-based features as a function of ρ. Similar curves for the other chord vocabularies can be found in Appendix B.2. We observe two results: The Baum–Welch-based features are superior to the Viterbi-based features for all values of ρ. In fact, even the worst value of ρ for the Baum–Welch algorithm performs better than the best for the Viterbi algorithm.

However, the assumption that a good chord recognition result leads to as good classification is obviously flawed. The Baum–Welch-based features achieve the maximal accuracy of 77.0% for $\rho = 0.3$. This is 5% higher than for the standard $\rho = 0.75$. We see that we can gain a 2.3% better performance than the combination of complexity and template features as proposed in [45]. However, we have to keep in mind, that we optimize this hyper-parameter on the *CrossEra* database and this particular feature setting. We do not use a separate cross validation, therefore the value of ρ is probably overfitted. We therefore continue to use $\rho = 0.75$ for the rest of the thesis since it lies in a relatively stable region. We assume that this holds also for other chord vocabularies. We also see that the inter-run deviation indicated by the error bars rises suddenly for $\rho \leq 0.15$. This indicates an unstable behavior of the features. We should therefore avoid Baum–Welch-based features with small values of ρ. Also, for large values of ρ, the Baum–Welch algorithm yields good features.

For the Viterbi algorithm, the optimal ρ is not far off the standard $\rho = 0.3$, therefore we continue to use it. We observe a drop of almost 10% for large values of ρ compared to standard ρ.

ρ	0.1	ρ_{std}	0.9	$\{0.1, \rho_{std}\}$	$\{\rho_{std}, 0.9\}$	$\{0.1, \rho_{std}, 0.9\}$	ρ_{opt}
Viterbi	69.9%	68.8%	62.1%	69.8%	67.5%	70.2%	70.0%
Baum–Welch	71.8%	72.6%	71.8%	75.7%	73.4%	73.9%	77.0%

Table 5.4: Results for combinations of features with different ρ.

This is probably due to the increasing sparsity of the feature vectors which we demonstrate in Figure 3.14. However, for $\rho = 0$, we still obtain an accuracy of 66.0%, which is quite high, considering the feature vector contains almost only noise-like transitions. Since the Viterbi-algorithm switches chords in each frame, those transitions which are not musically relevant predominate. In Figure 3.14 we see that most of those transitions are transitions between chords which have two notes in common. Interestingly, those transitions also seem to convey style information. Since Baum–Welch-based features also contain many of such transitions, this could be one reason for their good performance. To test this hypothesis, we delete all such transitions from the Baum–Welch-based feature vectors. However, we still obtain an accuracy of 72.4% for $\rho = 0.75$, which is only slightly worse than for the full feature vectors. We can therefore conclude that the high performance of the Baum–Welch-based features is not due to those "confusion transitions".

Features created with different values of ρ may have different semantic meaning. For small values of ρ, we measure the local behavior of chords, possibly taking non-chord notes into account. For large values of ρ, we measure chord changes on a larger scale, which can hint towards key changes. This information can possibly increase the performance of the classifier. We therefore combine features obtained with different values of ρ. By doing so, we hope to capture more information in the feature vector. On the other hand, we also increase the dimensionality of the feature vector, which also yields problems. For a good comparison between Viterbi- and Baum–Welch-based features we combine the features with the standard ρ and features with $\rho = 0.1$ and $\rho = 0.9$. We test combinations of the set $\{0.1, \rho_{std}, 0.9\}$, with $\rho_{std} = 0.3$ for the Viterbi-based features and $\rho_{std} = 0.75$ for the Baum–Welch-based features. Table 5.4 shows the results. As a reference we present results for all three individual values of ρ as well as for ρ_{opt}, the (overfitted) value for ρ that yields the best classification accuracy.

For the Baum–Welch-based features, we see that the results of the combined features are always higher than for the individual feature types. A combination of ρ_{std} with $\rho = 0.1$ works best, probably because $\rho_{std} = 0.75$ is already pretty high. However, we do not reach the maximal accuracy at $\rho_{opt} = 0.3$. For the Viterbi algorithm, a combination of all three values of ρ can exceed the maximum accuracy, though only by 0.2% and not enough to justify the increased effort. The other combinations do not lead to improved features. A reason for this may be that ρ_{std} is already close to ρ_{opt}.

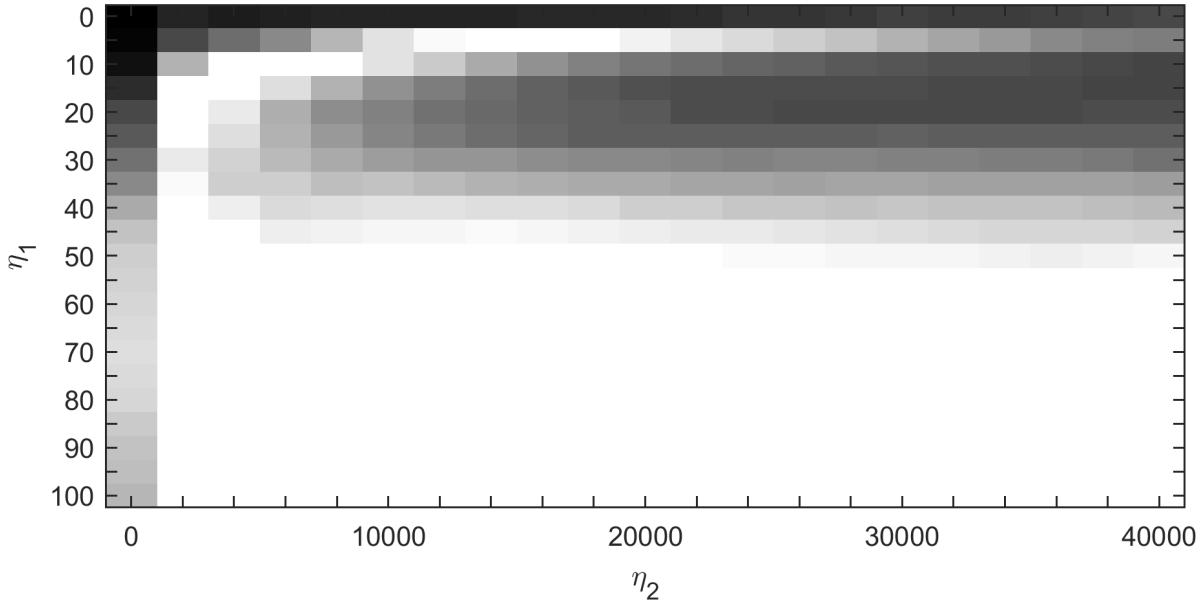


Figure 5.4: Classification accuracies for different rescaling parameters η_1, η_2.

5.6 Influence of Rescaling

In Section 3.2, we show that a non-linear rescaling improves the visual perception of local chord transitions features. For this purpose we propose a double-softmax rescaling of $\xi_n(i,j)$ which is able to enhance the "musically relevant" transitions, see Algorithm 3.2. In Section 3.4, we visually show the effect rescaling has on the piece-level chord transition features. We see that we are able to suppress transitions which probably were musically irrelevant. In this section, we test if there is also a qualitative improvement of the classification results when we apply double-softmax rescaling. For this purpose we generate features based on the *Triads* chord vocabulary using the Baum–Welch algorithm. Before feature aggregation we apply double-softmax rescaling on $\xi_n(i,j)$. We perform a grid search over the softmax parameters η_1 and η_2 with $\eta_1 \in \{0, 5, \ldots, 100\}$ and $\eta_2 \in \{0, 2000, \ldots, 40000\}$. Here, $\eta_{1/2} = 0$ denotes the case that we do not perform the respective softmax. A parameter set of $(\eta_1, \eta_2) = (0, 0)$ is therefore the case without any rescaling.

Figure 5.4 shows the result as an intensity plot. We see that the maximum accuracy lies at $(\eta_1, \eta_2) = (0, 0)$ with 72.6%. Even though rescaling with $(\eta_1, \eta_2) = (10, 10000)$ yields very good visual results for the Bach choral, we see that such a rescaling performs worse in a classification scenario with an accuracy of only 60.1%. We see that the performance gets better again for large η_2. For $(\eta_1, \eta_2) = (15, 40000)$ we obtain an accuracy of 69.0%. Even though the performance is still worse than without any rescaling, the existence of good accuracies for high η_2 is still interesting, so we further investigate this setting. For this purpose, we look at the confusion matrices for $(\eta_1, \eta_2) = (15, 40000)$ and compare it with the confusion matrix for $(\eta_1, \eta_2) = (0, 0)$

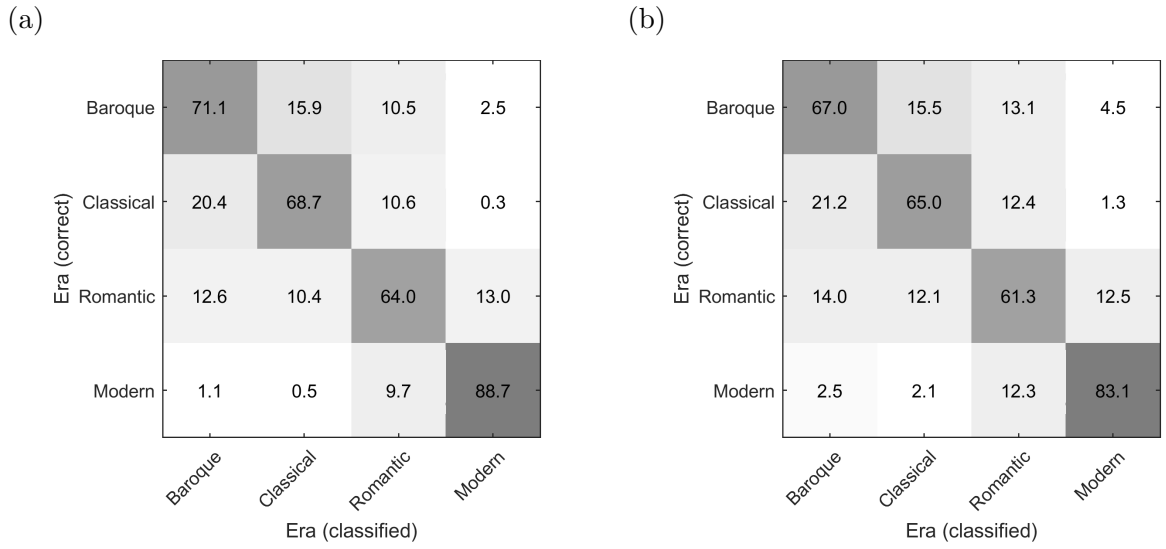


Figure 5.5: Confusion matrices for (η_1, η_2) equal to (a) $(0, 0)$ and (b) $(15, 40000)$.

in Figure 5.5. We see that all per-class accuracies are smaller for rescaling but the general behavior of era confusions does not change. We therefore conclude that the semantic information carried in the features is similar.

There are two possible reasons for the bad performance of the features vectors obtained with rescaling: It is possible that rescaling deletes important information. The other possibility is that the rescaling algorithm reacts too sensitive on the content and is therefore not able to perform equally for all recordings, such that the feature vectors differ too much, also within one class.

5.7 Reduction to Major and Minor Chord Types

When we compare the way of obtaining chord transition features in [45] with our method, we see several differences. In [45] Weiß performs chord estimation with a large chord vocabulary, before reducing the resulting chord labels to major and minor chord types. The idea behind this procedure is to only consider musically meaningful transitions, since forcing an augmented chord into either a major or minor triad is musically incorrect. To this end, we map only chords which were based on a major triad (also major seventh chords for example) to a major chord. We treat chords based on a minor triad accordingly. Chords which are neither based on a major nor a minor triads are ignored. In the following we define a chord as "based on a major or minor triad" if the three lowest notes form the respective triad.

Mapping	Flexible		Strict		None		
Orignal Vocabulary	*Triads*	*Sevenths*	*Triads*	*Sevenths*	*MajMin*	*Triads*	*Sevenths*
Viterbi	66.5%	62.4%	66.1%	61.3%	68.4%	68.8%	67.8%
Baum–Welch	70.5%	70.8%	67.9%	71.3%	69.1%	72.6%	74.6%

Table 5.5: Results of experiments with reduced chord vocabularies.

In our experiments with the *MajMin* vocabulary we only allow major or minor triads. Thus, we force every chord into the scheme of major and minor triads. We now want to reduce the *Triads* and *Sevenths* vocabulary to major and minor triads, in order to make the results comparable to the work in [45]. We can employ two strategies here. First, we map every chord to a major or minor triad according to the lowest third interval. This should resemble the strategy when we use the *MajMin* vocabulary. We call this strategy a "flexible mapping". The other strategy is taken from [45], where we tale only chords which are based on a major or minor triad into account. We discard all other chords and the respective transitions. We call this a "strict mapping". In Table 5.5 we show the results of experiments with different mappings. As a reference we give the accuracy obtained when we use the *MajMin* vocabulary already in the chord estimation stage, as well as the accuracies of the *Triads* and *Sevenths* chord vocabulary.

We see that the result with reduced vocabularies is always worse than any non-reduced vocabulary for the Viterbi algorithm. We observe that reduction from the *Sevenths* vocabulary is worse than from the *Triads* vocabulary. Note that the non-reduced *Sevenths* vocabulary also performs slightly worse than the *Triads* vocabulary. Furthermore, we see that the strict reduction performs worse than the flexible reduction, but only by a small difference. One explanation of why the flexible mapping performs worse than the *MajMin* vocabulary may be the following: For the flexible reduction we perform a mapping which is similar to the mapping the Viterbi algorithm implicitly performs when using the *MajMin* vocabulary. However, the Viterbi-algorithm optimizes the chord sequence in context. Performing the mapping after the algorithm ignores this context and may therefore lead to worse features. In general, we delete much information with the mapping. Since each frame contains only information about one transition, we delete the information contained in this frame completely for strict mapping.

Again the Baum–Welch-based features behave differently. Here, we see that the results improve by at least 1% compared to the *MajMin* vocabulary. An exception is the strict mapping of the *Triads* vocabulary. This may be due to the loss of information again. However, even for strict mapping we still conserve information from the frame, which describes major and minor chord types. For the Baum–Welch-based features a reduction from the *Sevenths* vocabulary performs better than from the *Triads* vocabulary. We need to keep in mind that features with a reduced chord vocabulary always show a worse performance than the features using the

Setting	No Filter	Composer Filter	Piano → Orchestra	Orchestra → Piano
Viterbi	66.4%	64.3%	67.1%	58.4%
Baum-Welch	77.4%	72.6%	69.4%	65.9%

Table 5.6: Style classification accuracies and inter-fold deviations (Dev.) for various features when different filters were applied.

respective non-reduced vocabulary. However, we show that we can improve the performance of major-minor-based feature types by mapping from a larger vocabulary.

5.8 Influence of the Album Effect

One result from [45] is that standard audio features such as audio spectral envelope, spectral flatness measure or mel frequency cepstral coefficients are not suitable for style classification, since those features mainly capture timbral information. Weiß [45] found that the features perform well without any constraints to the cross validation splits but failed with a composer filter. The conclusion was, that the classifier overfits due to recording artifacts of an album and uses these properties for classification. This effect is called "album effect". Employing a composer filter prevents that effect. Tonal features, on the other hand performed better with album filter.

In the previous sections, we already showed that our features work well also with a composer filter. In this section, we want to examine the effect of different cross validation strategies in detail. To this end, we compare the results in four different settings. First, we use no filter, which probably gives the best results since we can exploit recording artifacts and the personal style of the composer. Both properties simplify the separation of classes in *CrossEra*, but were excluded in the previous experiments on purpose. Second, we apply the composer filter as used in the previous sections. In the end, we train the classifier on *CrossEra-Piano* and test on *CrossEra-Orchestra* ("Piano → Orchestra") and vice versa ("Orchestra → Piano"). Since the timbral content of both instrumentation classes is different, those settings assure that we only use tonal information for classification. We summarize both settings under the name "instrumentation filter", but report both experiments separately. We test Viterbi-based and Baum–Welch-based features and show the results in Table 5.6.

As expected we see that the features perform better without filter than with composer filter. As previously discussed, the reason here is overfitting on non-musical information. We see that the drop of quality when we apply the composer filter is higher for Baum–Welch-based features than for Viterbi-based features. We assume that this is due to a larger amount of timbral information as a result from the soft-decision algorithm. For the instrumentation filter we look at the single

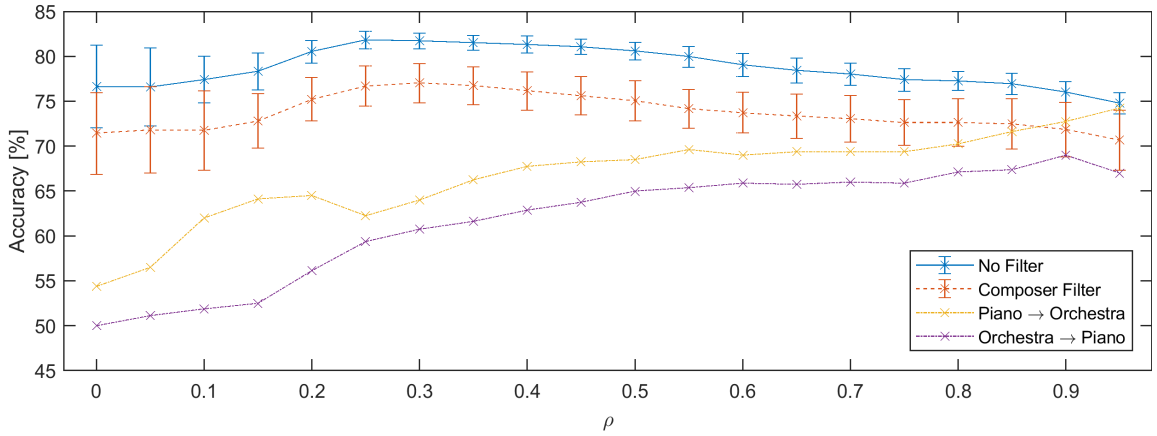
results of both possible setups. We see that training on piano recordings and testing on orchestra recordings yields better results than the reversed case. Weiß observes the same tendencies in [45]. Interestingly, for "Piano → Orchestra", the Viterbi-based features perform better than without filter, but the reversed case is roughly 10% worse. For the Baum–Welch-based features, both results are worse than the results with composer filter, but there is only a difference of 3.5% between both directions. This is a more stable behavior than the Viterbi-based features. We see that the instrumentation filter results are on average about 5% below the composer filter for Baum–Welch based features and less than 2% for Viterbi-based features. Those are better results than Weiß [45] obtains with template features, where accuracy dropped by roughly 10% compared to the composer filter.

We see that, even though the Viterbi-based features perform well under the instrumentation filter in average, the individual results show an unstable behavior. The Baum–Welch algorithm shows a different behavior. Here, the loss due to the instrumentation filter is higher, but the features perform comparably in both directions.

We observe another interesting difference when we perform the experiment for different self-transition probabilities ρ. In Figure 5.6, we show the accuracies of all four experiments as a function of ρ for both Baum–Welch-based features and Viterbi-based features. We observe that the difference between the curves for "no filter" and the composer filter is independent of the chosen value of ρ for both algorithms. For the Baum–Welch-based features, the performance of the instrumentation filter generally decreases with decreasing values of ρ. From the experiments on the Bach choral in Section 3.3, we know that there are more transitions present for smaller values of ρ, which points towards a larger amount of timbral information. This is one possible explanation for this result, since we assume that a larger amount of timbral information leads to worse generalization over different instrumentations. Interestingly, the curve for the composer filter does not decrease accordingly. We therefore conclude that the reason behind the loss of accuracy with composer filter is at least partially different to the reason behind the loss with instrumentation filter. Additionally, this plot justifies the choice of $\rho = 0.75$ as standard parameter for Baum–Welch-based features. The value of ρ which yields the highest result with the composer filter ($\rho = 0.3$) shows worse results with instrumentation filter. We therefore consider this feature type less robust against changes in instrumentation.

If we look at the Viterbi-based features, we see that the features behave unstable in an instrumentation filter scenario, in a sense of a large difference between the two cases, for all values of ρ. However, we see that—different to the Baum–Welch-based features—the loss due to the instrumentation filter does not depend on ρ. Since we connect this loss with the presence of timbral information, we conclude that the timbral content of Viterbi-based features is independent of the chosen value of ρ. An exception are features obtained with low values of ρ, where the accuracy

(a) Baum–Welch-based features



(b) Viterbi-based features

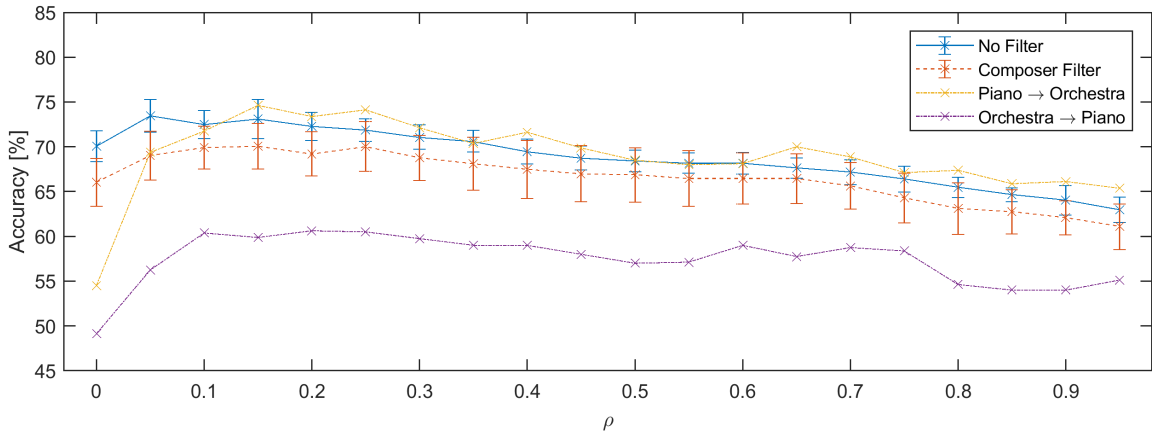


Figure 5.6: Classification accuracy of (a) Baum–Welch-based features and (b) Viterbi-based features for different ρ and different filters.

Era	Selected Composers
Baroque	J.S. Bach, Corelli
Classical	Boccherini, Haydn, Clementi
Romantic	Brahms, Grieg, R. Schumann, Brahms
Modern	Bartok, Berg, Schoenberg, Shostakovich

Table 5.7: Composers we chose for the test set of the cross-version experiment. The recordings of all remaining composers comprise the training set.

drops suddenly for the instrumentation filter. We showed in Section 3.3, that also Viterbi-based features capture much timbral content for small ρ, which is in line with our hypothesis.

In general, we can say that Baum–Welch-based features are influenced more by a composer or instrumentation filter. Nevertheless, Baum–Welch-based features still perform better than Viterbi-based features. Additionally, we observe an unstable behavior for the Viterbi-based features with instrumentation filter, since Viterbi-based features do not generalize well from orchestra data to piano data. We also found that the influence of the composer filter is weaker for chord transition features in general than for other tonal features proposed in [45].

5.9 Cross-Version Experiments

As a final experiment, we want to verify our classifier with cross-version experiments. For this purpose, we use the *CrossEraMirror* dataset. As described in Section 4.1, the database contains different recordings of a subset of the works in the *CrossEra* dataset. We did not use *CrossEraMirror* in previous experiment. We do not perform extensive research on *CrossEraMirror* to avoid presenting overfitted results. We use the *CrossEraMirror* database in two scenarios. First, we perform classification on one single split in training and test set, which was manually created to resemble the previous experiments as closely as possible. We use a composer filter and select composers for the test set, such that the test set contains roughly one third of the samples (compare Table 5.7). Afterwards, we discard all recordings that are not part of the *CrossEraMirror* database from the test set. We use the recordings from all the remaining composers for the training set. We stratify both sets using random deletion. Afterwards, we train a classifier on the training set by using recordings from *CrossEra* and test the classifier on recordings from both databases separately. That way we test, how well the features generalize over different versions. In this experiment, we expect that the result we obtain when we test on *CrossEra* data is better since recordings in *CrossEra* are from the same label as the training data. We test Viterbi-based and Baum–Welch-based features with all chord vocabularies. We give the results in Table 5.8. When evaluating this table we need to keep in mind, that we are dealing with only one classification experiment without cross validation.

	Viterbi			Baum-Welch		
	CrossEra	*CrossEraMirror*	Δ	*CrossEra*	*CrossEraMirror*	Δ
MajMin	71.2%	72.7%	-1.5%	70.6%	69.0%	1.5%
Triads	67.9%	68.5%	-0.6%	72.7%	70.6%	2.1%
Seventh	67.1%	65.2%	1.9%	75.0%	70.8%	4.2%
Intervals	61.3%	61.2%	0.2%	71.3%	69.8%	1.5%

Table 5.8: Results of the first experiment on *CrossEraMirror*. We use one manually created split. We always train the model on *CrossEra* data and test on both datasets.

Surprisingly, the Viterbi-algorithm yields sometimes better results on *CrossEraMirror* than on *CrossEra*. Since the deviations are comparatively small and go in both directions, it is possible that the deviation is within the normal inter-fold deviation. In general, we see that Viterbi-based features generalize well over different versions. When we average over all types of Viterbi-based features, there is no difference between both databases.

The Baum–Welch-based features generalize worse over different versions. If we use the *CrossEraMirror* database for testing, we see that the results are consistently worse than for *CrossEra*. We observe a trend that the difference increases with larger chord vocabularies. In contrast, the *Intervals* vocabulary shows a lower deviation again.

Since both databases differ mainly in timbral content—for example, *CrossEraMirror* contains recordings of an harpsichord, which was excluded in *CrossEra*—this experiment supports the hypothesis, that Baum–Welch-based features contain more timbral content. We have to keep in mind that—except for the *MajMin* dataset—the Baum–Welch-based features still perform better in terms of achieved accuracy.

For the second experiment, we use the *CrossEraMirror* database to simulate a "real" album filter. Therefore, again, we train a model on data from *CrossEra* and then test it on *CrossEraMirror*. We reduced both datasets to the works contained in *CrossEraMirror*. This way we can perform a cross validation on both sets. We do not use any filter here, but since *CrossEra* and *CrossEraMirror* contain no common album, the cross-version experiment contains an implicit album filter. We show the results in Table 5.9. We see that most feature types have a similar behavior and loose between 2% and 3% accuracy when we test them on *CrossEraMirror*. Reasons can include the album effect or differences in instrumentation (grand piano in *CrossEra* and harpsichord in *CrossEraMirror*). An exception are the Viterbi-based features on the *MajMin* vocabulary and the Baum–Welch-based features on the *Triads* vocabulary, which perform better. We have no explanation for these outliers. We cannot compare the results directly to the results from the previous section, where we already showed experiments on *CrossEra* without any filter, since the used datasets differ in various points, such as size and the distribution of the composers (also compare Table 4.1). That way, we are dealing with a different task here.

	Viterbi			Baum-Welch		
	CrossEra	*CrossEraMirror*	Δ	*CrossEra*	*CrossEraMirror*	Δ
MajMin	72.5%	72.0%	0.5%	74.6%	72.4%	2.2%
Triads	73.7%	71.3%	2.4%	79.7%	78.2%	1.5%
Seventh	73.1%	70.1%	3.0%	82.3%	79.6%	2.7%
Intervals	68.8%	66.3%	2.6%	80.6%	78.2%	2.4%

Table 5.9: Results of the second experiment performed on *CrossEraMirror*, which simulates an album filter. We always train the model on *CrossEra* data but test on either *CrossEra* or *CrossEraMirror*.

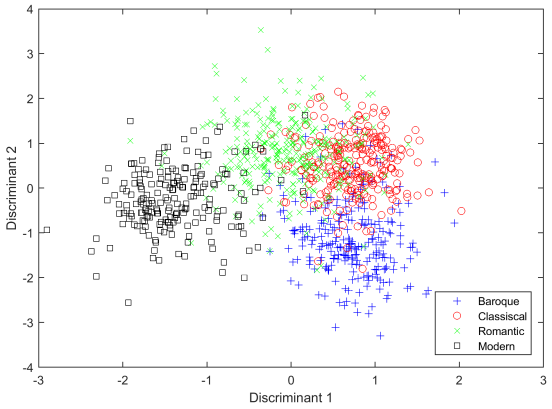
We also perform the same experiment using the template features from [45]. We obtain an accuracy of 76.0% when we test those features on *CrossEraMirror* as opposed to an accuracy of 82.8% on *CrossEra*. We see that chord transition features achieve better generalization than template features. This is in line with the results in Section 5.8, where we examine generalization over instrumentation.

To show the relationship between the features of both datasets, we use a scatter plot of the reduced features after LDA in Figure 5.7. For this plot, we use Baum–Welch-based features on the *Triads* vocabulary. We first look at the left column of plots. Here, we see the features from the *CrossEra* dataset. To visualize the 3D-data obtained from LDA, each plot shows a different combination of coordinates. Note that the classes are usually better separable than suggested by a single plot. One plot shows only two coordinates, but a classifier can use all three coordinates for classification. For the plots, we fitted LDA on the entire *CrossEra* dataset. The classes are color-coded. The plots confirm our conclusions drawn from the confusion matrices in Section 5.4. The classes Baroque, Classical, and Romantic are distinct but have overlapping regions. Those overlaps are present in all plots. We see that the Modern class is clearly separable from Baroque and Classical by the first discriminant but still has a few overlaps with Romantic. This is also consistent with the previous results and the findings in [45].

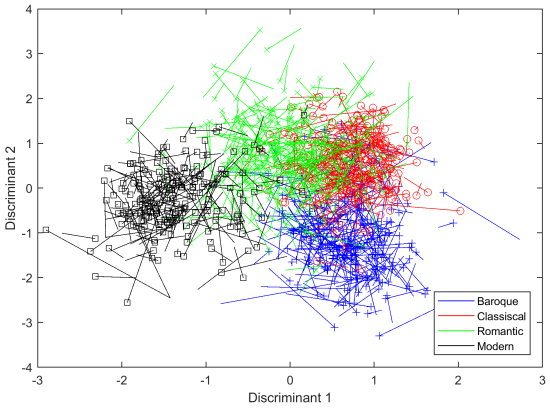
On the right column we see an extension of the previous plots. Here, we show the same features from *CrossEra* but we show the distance between the musically corresponding instances from *CrossEraMirror* as a line connecting both instances. For orientation purposes, we mark the instance from *CrossEra*. We see that the distance between the features can get rather large. However the lines rarely cross the boundaries between the classes. This indicates a stable work-level behavior in classification experiments.

In Figure 5.8 we show the same plot for features based on the *Sevenths* chord vocabulary. Again, we use Baum–Welch-based features. We only display the projection on a single plane for both datasets in one plot. We see that the classes are better separable than for the *Triads* dataset. This is in line with the higher accuracies we obtain with the *Sevenths* vocabulary. However, we

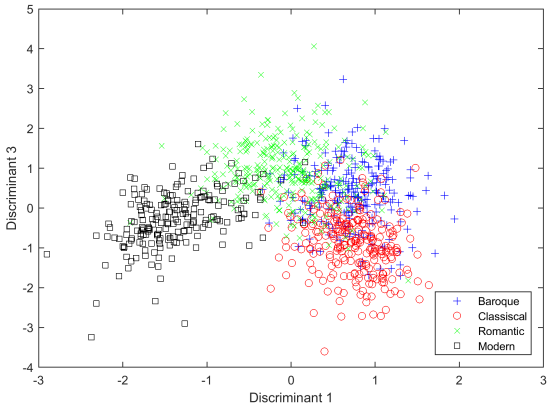
(a) *CrossEra*



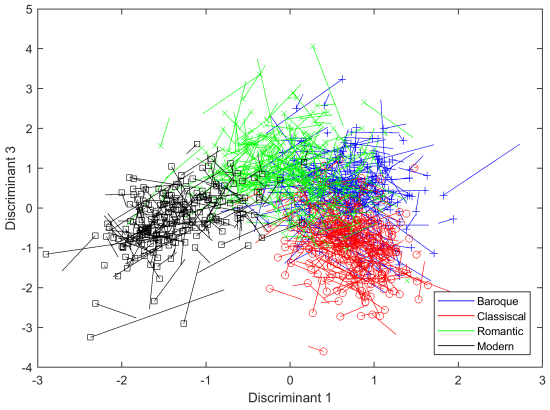
(b) *CrossEra*—*CrossEraMirror*



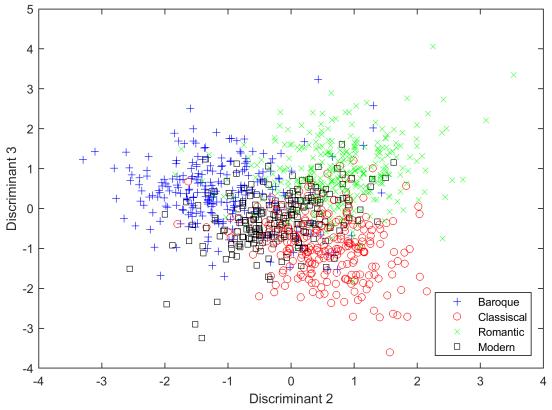
(c)



(d)



(e)



(f)

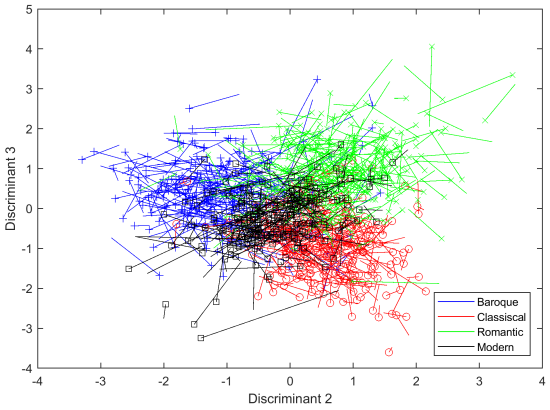


Figure 5.7: Scatter plot of all instances from the *CrossEra* database. In the left column, each point corresponds to one recording in the *CrossEra* database. In the right column, each line connects an instance vector from *CrossEra* with the musically corresponding instance from *CrossEraMirror*. Each row shows the projection of the 3D-data on a different coordinate plane. We use features computed with the Baum–Welch algorithm and the *Triads* chord vocabulary.

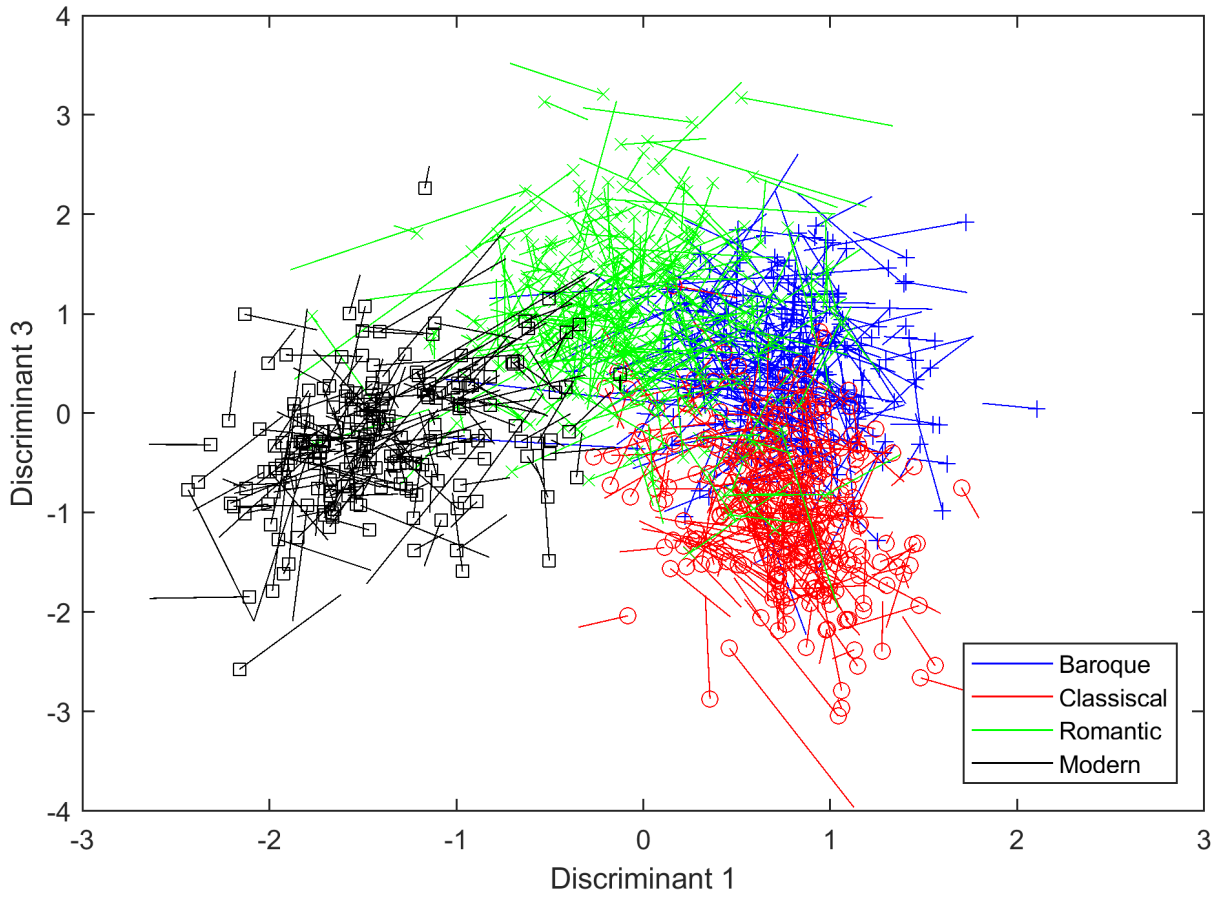


Figure 5.8: Scatter plot showing the features of *CrossEra* connected with the features of *CrossEraMirror*. We compute the features with the Baum–Welch algorithm and the *Sevenths* chord vocabulary.

see several lines crossing the class boundaries. Most prominently, we see many features "moving" from Modern to Romantic and from Romantic to Classical. That visualizes the difference between the results of both datasets, which is also in line with the loss of accuracy we observe in Table 5.8 for the *Sevenths* vocabulary. We see that this kind of plot is suitable for showing the difference between the two datasets.

In this section, we showed that chord transition features generally perform well in cross-version experiments. We found differences between Baum–Welch-based features and Viterbi-based features. The high-level Viterbi-based features generalize better over different versions. We assume, this is due to only few timbral information. The results we presented here are in line with the results from the previous sections.

Chapter 6

Conclusions

In this thesis, we introduced a novel kind of mid-level chord transition features. Chord transition features are useful for describing the harmonic progression of music recordings and contain information about the frequency of different kinds of chord transitions. Chord transition features described in previous work are high-level features, which directly describe properties of an estimated chord sequence. This approach is based on chord recognition, which is usually done with Hidden Markov Models and the Viterbi algorithm. In this thesis, we propose mid-level features. We do not estimate a chord sequence from which we derive the transition features, but rather compute transition probabilities based on the probabilistic Hidden Markov Model. With this kind of features, we are able to take ambiguities in the chord estimation process into account.

To compute local mid-level chord transition features, we made use of the Baum–Welch algorithm. This algorithm was initially proposed to train a Hidden Markov Model. We derived state transition probabilities from internal variables of the Baum–Welch algorithm. Extracting these variables, we were able to perform a soft-decision chord transition estimation. For visualizing the local mid-level chord transitions, we proposed a non-linear rescaling technique. In a chord recognition scenario, we showed that this approach is able to display ambiguities in the recognition process. We used these ambiguities to gain more musical knowledge than from a traditional hard-decision chord recognition approach.

We also showed that we can use chord transitions to characterize a whole recording. We obtain these piece-level chord transition features by averaging the chord-transition probabilities over the entire sequence. We test these aggregated features in a style classification scenario, in which we distinguish between the four musical periods Baroque, Classical, Romantic, and Modern within Western classical music.

In the classification experiments, the proposed Baum–Welch-based chord transition features showed consistently better classification results than equivalent high-level Viterbi-based features. Baum–Welch based features benefit from using a larger chord vocabulary, while the efficiency of Viterbi-based features decreases with larger chord vocabularies. Even though chord transitions are only one aspect to distinguish musical styles, our mid-level features show a performance comparable to other state-of-the-art tonal features. We showed that a combination of the mid-level chord transition features with those state-of-the-art tonal features performed better than both individual feature types alone. We performed all experiments using a composer filter, to prevent overfitting to non-musical properties known as the "album effect". We also tested the features without composer filter and found that the album effect has less impact than on other mid-level tonal features. We also performed experiments in which we trained the classifier on piano pieces and tested on orchestra pieces and vice versa. Here, we showed that the Baum–Welch-based features are in general robust to different instrumentation. In contrast, Viterbi-based features showed an unstable behavior.

There are several interesting topics concerning mid-level chord transition features, which could be examined in future work. We found that the self-transition probability is an important hyper-parameter of the Hidden Markov Model we used to generate our features. We observed effects on the classification accuracies as well the features ability to generalize over different instrumentations. In future work, one could further investigate the internal mechanisms leading to this behavior and search for optimal values in a separate cross validation.

In this thesis we limited ourself to a template emission model. Cho and Bello [9] found that a Gaussian emission model is generally more robust against changes in the self-transition probability. We showed that a Gaussian model allows chord vocabularies containing chords with different numbers of notes. Researching the applications of a Gaussian emission model to generate mid-level chord transition features is therefore promising interesting results.

The usage of Hidden Markov Models gives us another interesting possibility for a different classifier. Instead of creating features with the Baum–Welch algorithm, we could use this algorithm to train one HMM for each class. Afterwards, we could classify a recording by evaluating the probability of the respective chroma sequence within each model and choosing the class whose model yields the highest probability. A similar approach is typically used for speech recognition tasks, such as isolated word recognition [33]. However, first experiments did not show promising results for style classification. In future work one could further investigate this method.

In this thesis we showed the high performance or our features in one specific scenario: classifying recordings of Western classical music into four sub-styles. In future work we could examine the features in other scenarios, such as in jazz or rock music. Also, changing the classification objective, such as classifying different composers or looking for applications in work identification is possible.

In summary, this thesis shows that it is possible to combine the benefits of chord transition features with the concept of mid-level, soft-decision features. The resulting features are robust against the album effect, generalize well across instrumentations, and perform better than comparable high-level chord transition features, based on hard-decision chord estimation.

Appendix A

Hidden Markov Models

In this chapter, we derive the equations for the forward and backward algorithm and the calculation of the local state transition probabilities $\xi_n(i,j)$, which were introduced in Section 3.2. We use the notations and definitions introduced there.

A.1 Forward Algorithm

First, we recall the definition of the forward variable $v_n(i)$ (Eq. 3.13):

$$v_n(i) = P(\boldsymbol{o}_1 \boldsymbol{o}_2 \dots \boldsymbol{o}_n, s_n = \alpha_i | \Theta) \tag{A.1}$$

For $n = 1$ we can write:

$$v_1(i) = P(\boldsymbol{o}_1, s_1 = \alpha_i | \Theta) = P(s_1 = \alpha_i | \Theta)\, P(\boldsymbol{o}_1 | s_1 = \alpha_i, \Theta) = c_i b_i(\boldsymbol{o}_1), \tag{A.2}$$

using the chain rule. That way we arrive at the initialization step. To derive the iteration step, we need to use the Markov property. It states, that the current state only depends on the previous state. Let O_a^b with $1 \leq a < b \leq N$ denote the subsequence $[\boldsymbol{o}_a \boldsymbol{o}_{a+1} \dots \boldsymbol{o}_b]$. We write:

$$\begin{aligned} v_n(j) &= P\big(O_1^{n-1} \boldsymbol{o}_n, s_n = \alpha_j | \Theta\big) = \\ &= P\big(O_1^{n-1}, s_n = \alpha_j | \Theta\big)\, P\big(\boldsymbol{o}_n | O_1^{n-1}, s_n = \alpha_j, \Theta\big) \end{aligned} \tag{A.3}$$

Since $\boldsymbol{o}_n$ only depends on s_n and not directly on the previous observations (the observations are not statistically independent but all information passes through s_n, therefore all other conditions are redundant). We can write

$$P\big(\boldsymbol{o}_n | O_1^{n-1}, s_n = \alpha_j, \Theta\big) = P(\boldsymbol{o}_n | s_n = \alpha_j, \Theta) = b_j(\boldsymbol{o}_n) \tag{A.4}$$

We can rewrite the other part of Eq. A.3 by reverse marginalization:

$$P\big(O_1^{n-1}, s_n = \alpha_j | \Theta\big) = \sum_{i=1}^{I} P\big(O_1^{n-1}, s_n = \alpha_j, s_{n-1} = \alpha_i | \Theta\big)\,. \tag{A.5}$$

We now see that

$$\begin{aligned} P\big(O_1^{n-1}, s_n = \alpha_j, s_{n-1} = \alpha_i | \Theta\big) &= P\big(s_n = \alpha_j | O_1^{n-1}, s_{n-1} = \alpha_i, \Theta\big)\, P\big(O_1^{n-1}, s_{n-1} = \alpha_i | \Theta\big) = \\ &= P(s_n = \alpha_j | s_{n-1} = \alpha_i, \Theta)\, P\big(O_1^{n-1}, s_{n-1} = \alpha_i | \Theta\big) = \\ &= a_{ij} v_{n-1}\,(i)\,. \end{aligned} \tag{A.6}$$

In the second line we use the Markov property again to eliminate a condition. By combining Eqs. A.3–A.6, we obtain the known iteration step formula.

A.2 Backward Algorithm

We now derive the equations for the backward algorithm. Here, we have the problem that according to the definition

$$w_n\,(i) = P(\boldsymbol{o}_{n+1}\boldsymbol{o}_{n+2}\ldots\boldsymbol{o}_N | s_n = \alpha_i, \Theta) \tag{A.7}$$

the initial variable $w_N\,(i)$ is not defined. We therefore first derive the iteration step and then show, that $w_N\,(i) = 1\ \forall i$ yields the desired result. We start be rewriting the definition with the shorthand notation $O_a^b = [\boldsymbol{o}_a\boldsymbol{o}_{a+1}\ldots\boldsymbol{o}_b]$ with $1 \leq a < b \leq N$ and applying a reversed marginalization:

$$\begin{aligned} w_n\,(i) &= P(\boldsymbol{o}_{n+1}\boldsymbol{o}_{n+2}\ldots\boldsymbol{o}_N | s_n = \alpha_i, \Theta) \\ &= P\big(\boldsymbol{o}_{n+1} O_{n+2}^N | s_n = \alpha_i, \Theta\big) \\ &= \sum_{j=1}^{I} P\big(\boldsymbol{o}_{n+1} O_{n+2}^N, s_{n+1} = \alpha_j | s_n = \alpha_i, \Theta\big)\,. \end{aligned} \tag{A.8}$$

We can further expand the summands by applying the chain rule:

$$\begin{aligned} &P\big(\boldsymbol{o}_{n+1} O_{n+2}^N, s_{n+1} = \alpha_j | s_n = \alpha_i, \Theta\big) = \\ =&P(s_{n+1} = \alpha_j | s_n = \alpha_i, \Theta)\, P\big(\boldsymbol{o}_{n+1} O_{n+2}^N | s_{n+1} = \alpha_j, s_n = \alpha_i, \Theta\big) = \\ =&a_{ij} \cdot P\big(\boldsymbol{o}_{n+1} O_{n+2}^N | s_{n+1} = \alpha_j, \Theta\big)\,. \end{aligned} \tag{A.9}$$

In the last step we use the definition of the transition probability and the fact that the future output only depends on the previous state through s_{n+1}, which follows from the Markov property. We continue with

$$\begin{aligned} & P\big(\boldsymbol{o}_{n+1}O_{n+2}^{N}|s_{n+1}=\alpha_j,\Theta\big) = \\ =&P(\boldsymbol{o}_{n+1}|s_{n+1}=\alpha_j,\Theta)\,P\big(O_{n+2}^{N}|\boldsymbol{o}_{n+1},s_{n+1}=\alpha_j,\Theta\big) = \\ =&P(\boldsymbol{o}_{n+1}|s_{n+1}=\alpha_j,\Theta)\,P\big(O_{n+2}^{N}|s_{n+1}=\alpha_j,\Theta\big) = \\ =&b_j(\boldsymbol{o}_{n+1})w_{n+1}\,(j) \end{aligned} \tag{A.10}$$

Combining all above equations we arrive at the iteration step. We can explicitly compute $w_{N-1}\,(i)$:

$$\begin{aligned} w_{N-1}\,(i) &= P(\boldsymbol{o}_N|s_{N-1}=\alpha_i,\Theta) = \\ &= \sum_{j=1}^{I} P(\boldsymbol{o}_N, s_N=\alpha_j|s_{N-1}=\alpha_i,\Theta) = \\ &= \sum_{j=1}^{I} P(s_N=\alpha_j|s_{N-1}=\alpha_i,\Theta)\,P(\boldsymbol{o}_N|s_N=\alpha_j,s_{N-1}=\alpha_i,\Theta) = \\ &= \sum_{j=1}^{I} a_{ij}b_j(\boldsymbol{o}_N)\cdot 1. \end{aligned} \tag{A.11}$$

The calculation steps follow the same scheme as the previous derivation. We see that this expression is equivalent to the iteration step, if we set $w_N\,(i)=1\ \forall i$. Therefore we use this as initial value for the iteration.

A.3 Local State Transition Probabilities

To prove Equation 3.15 we first rewrite the conditional probability of Eq. 3.12 using Bayes' theorem:

$$\begin{aligned} \xi_n\,(i,j) &= P(s_n=\alpha_i, s_{n+1}=\alpha_j|O,\Theta) = \\ &= \frac{P(s_n=\alpha_i, s_{n+1}=\alpha_j, O|\Theta)}{P(O|\Theta)} \end{aligned} \tag{A.12}$$

In the following we only consider the enumerator. Let O_a^b with $1 \leq a < b \leq N$ denote the subsequence $[\boldsymbol{o}_a \boldsymbol{o}_{a+1} \dots \boldsymbol{o}_b]$.

$$\begin{aligned} &P(s_n{=}\alpha_i, s_{n+1}{=}\alpha_j, O|\Theta) = \\ {=}&P\big(s_n{=}\alpha_i, s_{n+1}{=}\alpha_j, O_1^n O_{n+1}^N|\Theta\big) = \\ {=}&P(O_1^n, s_n{=}\alpha_i, s_{n+1}{=}\alpha_j|\Theta) \cdot P\big(O_{n+1}^N|O_1^n, s_n{=}\alpha_i, s_{n+1}{=}\alpha_j, \Theta\big) \end{aligned} \tag{A.13}$$

Using the Markov property, we see that any observation at time $n+1$ or later only depends on previous states and observations through the state s_{n+1}. Of course they are not statistically independent but all information passes through s_{n+1}, therefore all other conditions are redundant. With this consideration we write:

$$\begin{aligned} &P\big(O_{n+1}^N|O_1^n, s_n{=}\alpha_i, s_{n+1}{=}\alpha_j, \Theta\big) = \\ {=}&P\big(O_{n+1}^N|s_{n+1}{=}\alpha_j, \Theta\big) = \\ {=}&P\big(\boldsymbol{o}_{n+1} O_{n+2}^N|s_{n+1}{=}\alpha_j, \Theta\big) = \\ {=}&P(\boldsymbol{o}_{n+1}|s_{n+1}{=}\alpha_j, \Theta) \cdot P\big(O_{n+2}^N|\boldsymbol{o}_{n+1}, s_{n+1}{=}\alpha_j, \Theta\big) \\ {=}&P(\boldsymbol{o}_{n+1}|s_{n+1}{=}\alpha_j, \Theta) \cdot P\big(O_{n+2}^N|s_{n+1}{=}\alpha_j, \Theta\big) \end{aligned} \tag{A.14}$$

In the last step we can eliminate one condition by similar considerations as above. Using the definitions of the backward variable (Eq. 3.14) and emission probability (Eq. 2.16) we see:

$$\begin{aligned} &P(\boldsymbol{o}_{n+1}|s_{n+1}{=}\alpha_j, \Theta) \cdot P\big(O_{n+2}^N|s_{n+1}{=}\alpha_j, \Theta\big) = \\ {=}&P(\boldsymbol{o}_{n+1}|s_{n+1}{=}\alpha_j, \Theta) \cdot P(\boldsymbol{o}_{n+2} \dots \boldsymbol{o}_N|s_{n+1}{=}\alpha_j, \Theta) = \\ {=}&b_j(\boldsymbol{o}_{n+1}) \cdot w_{n+1}(j) \end{aligned} \tag{A.15}$$

We now further analyse $P(O_1^n, s_n{=}\alpha_i, s_{n+1}{=}\alpha_j|\Theta)$ from Eq. A.13

$$\begin{aligned} &P(O_1^n, s_n{=}\alpha_i, s_{n+1}{=}\alpha_j|\Theta) = \\ {=}&P(O_1^n, s_n{=}\alpha_i|\Theta) \cdot P(s_{n+1}{=}\alpha_j|O_1^n, s_n{=}\alpha_i, \Theta) = \\ {=}&P(O_1^n, s_n{=}\alpha_i|\Theta) \cdot P(s_{n+1}{=}\alpha_j|s_n{=}\alpha_i, \Theta) \end{aligned} \tag{A.16}$$

Again one condition could be eliminated by the Markov property, this time by exploiting that a state is connected to previous observations only via the previous state. The resulting expression can be directly rewritten as

$$P(O_1^n, s_n{=}\alpha_i|\Theta) \cdot P(s_{n+1}{=}\alpha_j|s_n{=}\alpha_i, \Theta) = v_n(i) \cdot a_{ij}. \tag{A.17}$$

Combining Eqs. A.12 – A.17 we arrive at the expression in Eq. 3.15.

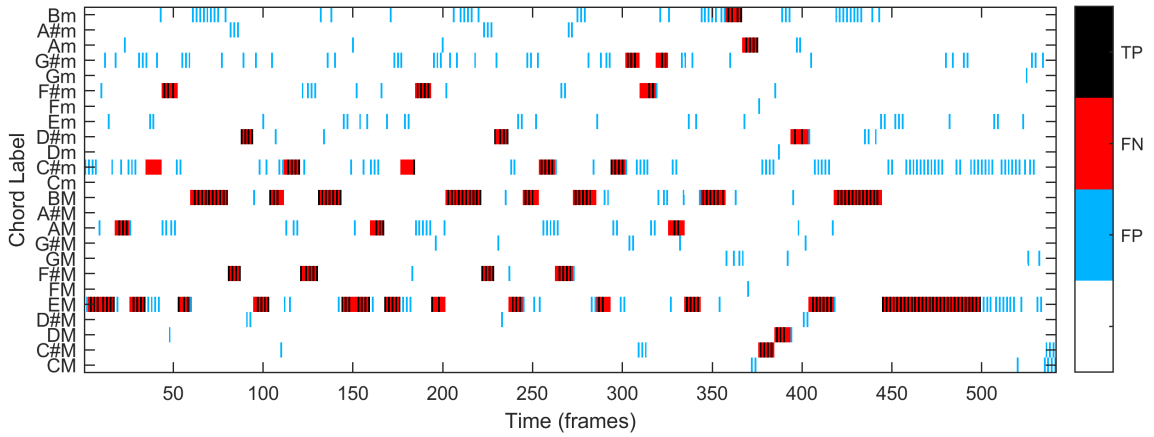
Appendix B

Further Results

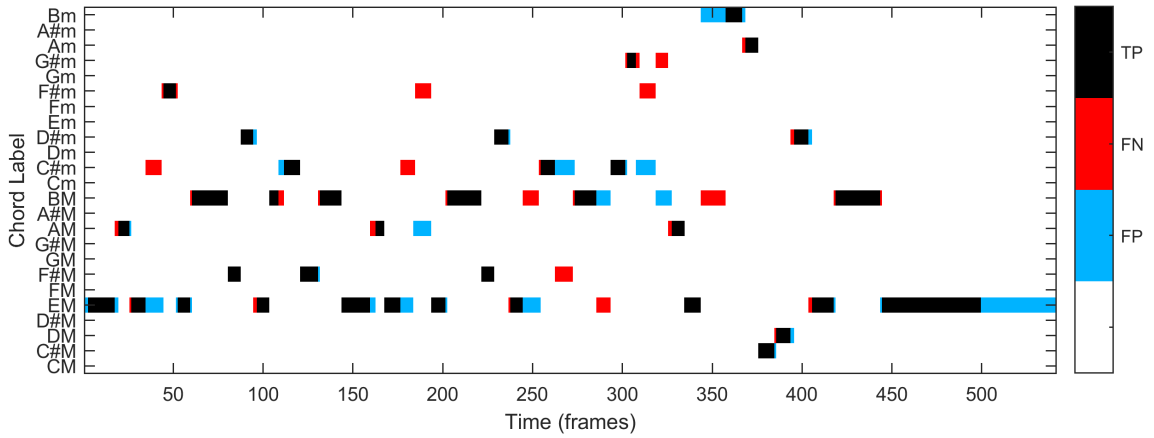
B.1 Influence of the Self-Transition Probability on Chord Recognition

In this section, we show plots to demonstrate the chord recognition performance of the Viterbi algorithm and the Baum–Welch algorithm for different values of ρ. In Figure B.1, we show the results of the Viterbi algorithm and in Figure B.2, we show the soft-decision results obtained with the Baum–Welch-algorithm. Here, we use a single softmax $\eta = 10$ on the state probabilities to visualize the probabilities better. As in Figure 3.7 we choose the parameter ρ from the set $\rho \in \{0.025, 0.02, 0.5\}$.

(a) $\rho = 0.025$



(b) $\rho = 0.2$



(c) $\rho = 0.5$

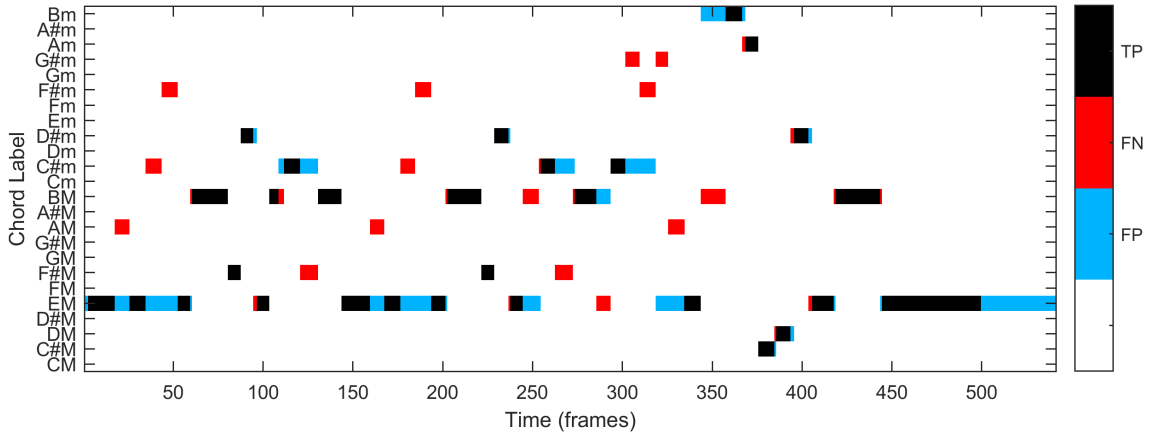
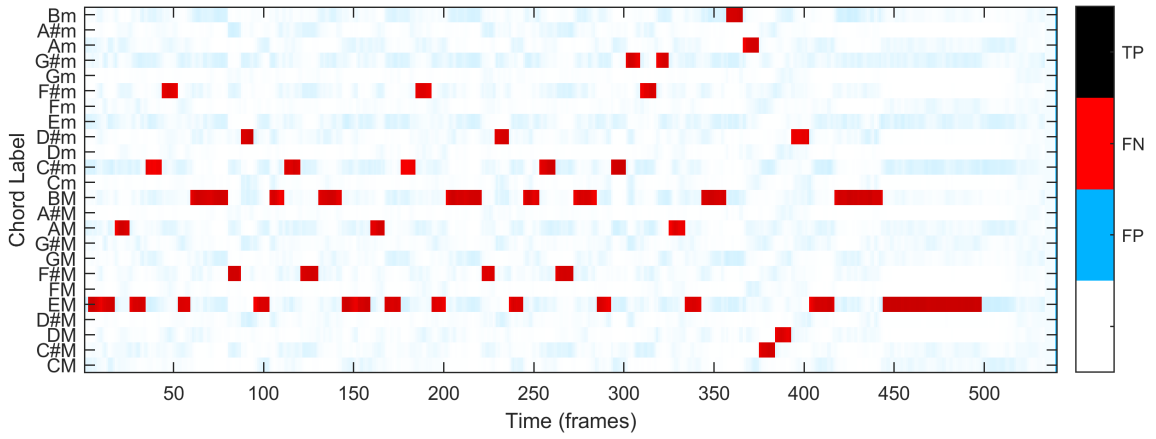
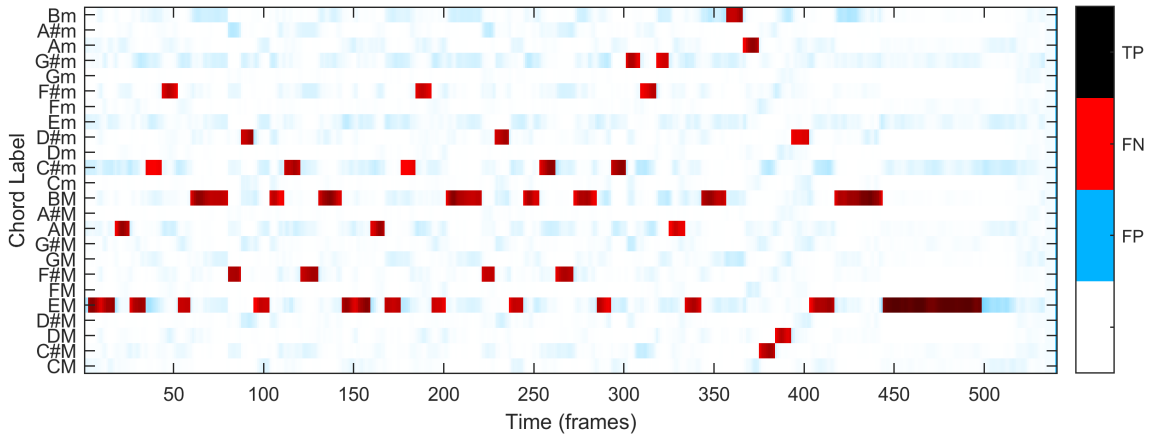


Figure B.1: Viterbi-based chord recognition of the Bach example for different values of ρ.

(a) $\rho = 0.025$



(b) $\rho = 0.2$



(c) $\rho = 0.5$

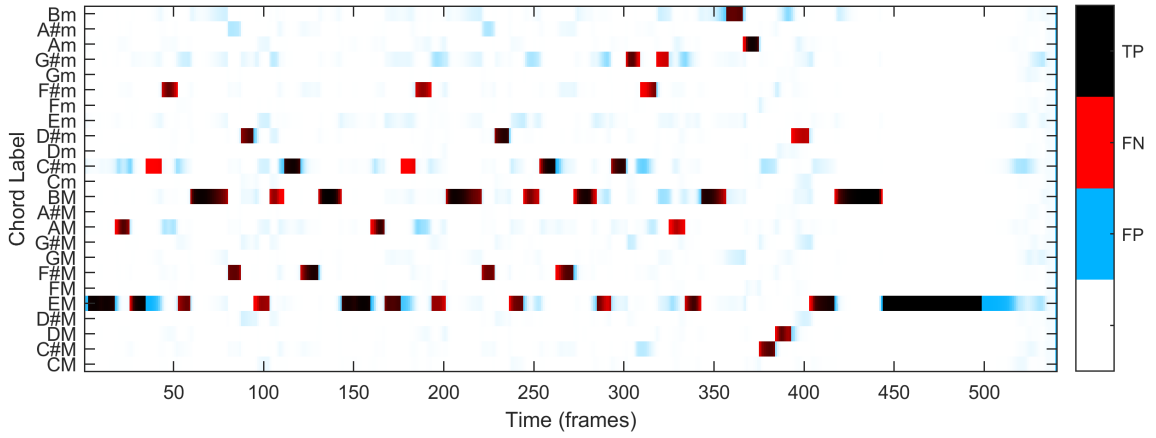
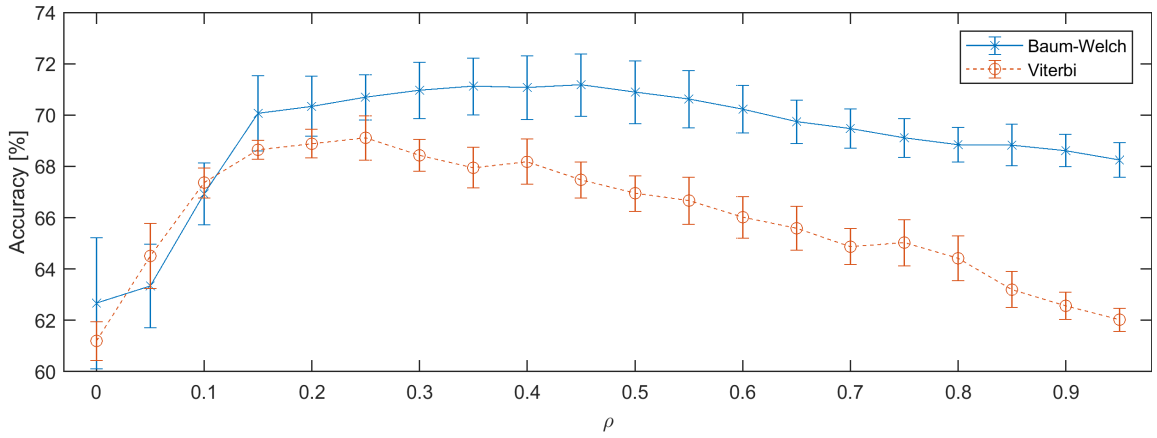


Figure B.2: Baum–Welch-based chord recognition of the Bach example for different values of ρ.

B.2 Influence of the Self-Transition Probability on Style Classification

In this section, we show several plots which show the influence ρ has on the classification accuracy. Figure B.3 shows the results for the *MajMin* and *Triads* vocabularies, and Figure B.4 shows the results for the *Sevenths* and *Intervals* vocabularies. The plots support the results obtained in Section 5.6.

(a) *MajMin*



(b) *Triads*

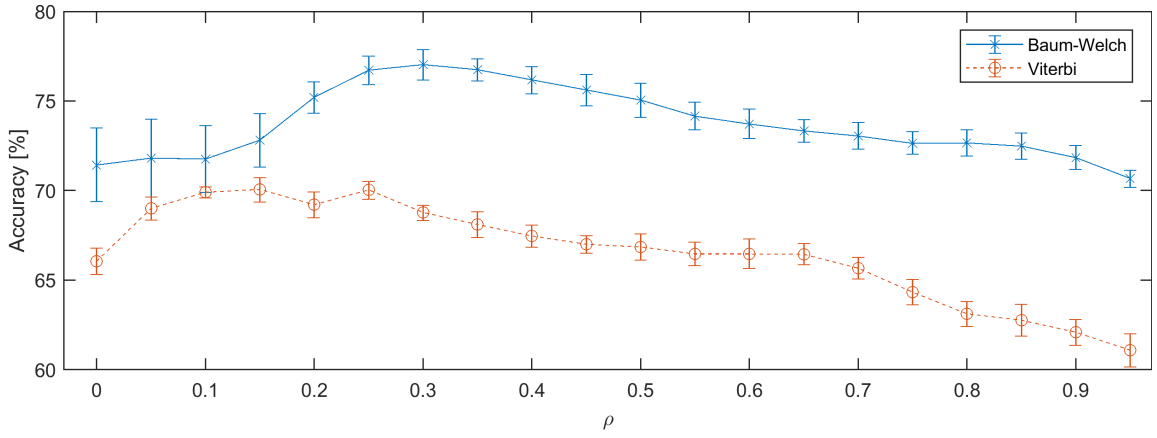
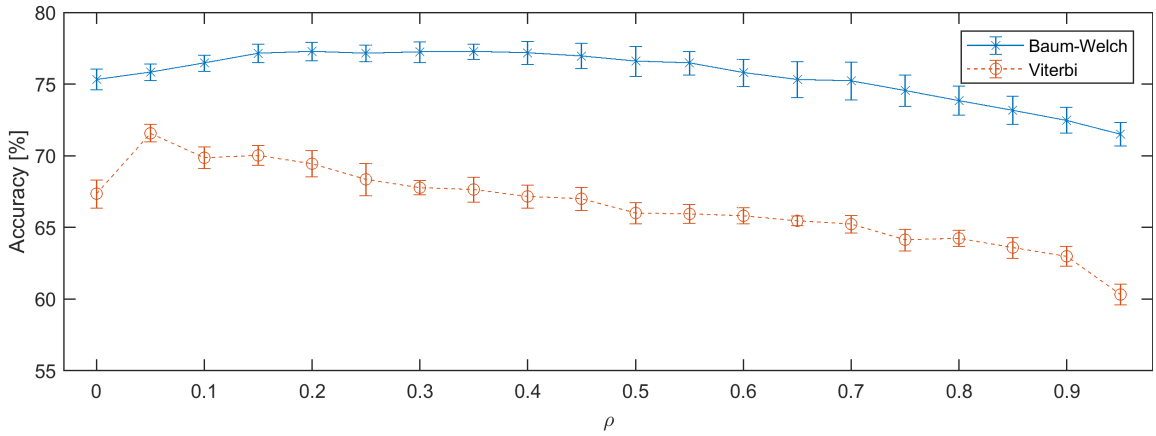


Figure B.3: Classification accuracy as a function of ρ for the *MajMin* and *Triads* vocabularies.

(a) *Sevenths*



(b) *Intervals*

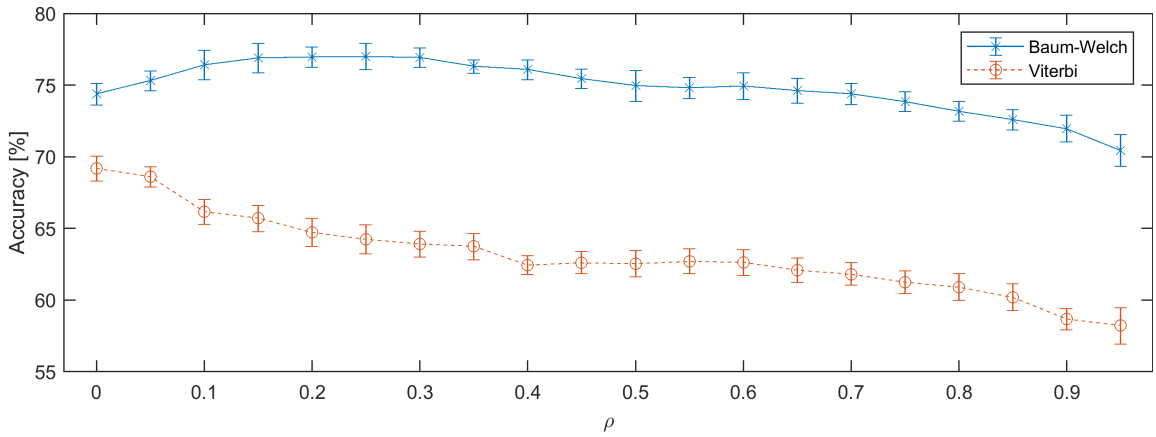


Figure B.4: Classification accuracy as a function of ρ for the *Sevenths* and *Intervals* vocabularies.

B.3 Further Classification Results

This section contains further results of the style classification experiments. Table B.1 gives results for all chord vocabularies, combinations with tonal feature types [45] and as a reference the individual tonal features types alone.

Feature Type / Chord Vocabulary	Viterbi			Baum–Welch		
	Full	*Piano*	*Orchestra*	*Full*	*Piano*	*Orchestra*
Chord Vocabularies						
MajMin	68.4%	60.4%	75.0%	69.1%	62.8%	75.7%
Triads	68.8%	58.3%	72.8%	72.6%	64.4%	76.2%
Sevenths	67.8%	61.6%	72.7%	74.6%	68.1%	78.6%
Intervals	63.9%	62.4%	66.1%	73.9%	67.7%	77.6%
Combination With Template Features						
MajMin	76.1%	73.2%	81.5%	76.7%	73.5%	82.0%
Triads	74.4%	65.6%	78.2%	76.0%	72.4%	82.3%
Sevenths	69.1%	63.9%	77.2%	76.9%	72.0%	82.7%
Intervals	68.6%	65.7%	74.9%	75.0%	71.3%	81.3%
Combination With Complexity Features						
MajMin	73.5%	71.4%	78.7%	74.9%	72.7%	80.5%
Triads	73.4%	60.3%	77.7%	75.9%	69.0%	81.5%
Sevenths	68.0%	62.1%	74.3%	75.3%	69.9%	82.6%
Intervals	66.5%	63.7%	70.9%	75.4%	69.9%	82.3%
Combination With Complexity & Template Features						
MajMin	77.5%	71.7%	81.8%	78.0%	72.5%	82.6%
Triads	74.0%	67.1%	79.5%	76.1%	72.8%	82.9%
Sevenths	69.9%	65.4%	78.0%	77.6%	73.0%	83.6%
Intervals	69.8%	66.9%	75.6%	76.4%	72.1%	82.9%
No Composer Filter						
MajMin	70.9%	64.0%	76.4%	73.5%	69.5%	79.4%
Triads	71.1%	64.3%	75.3%	77.4%	74.2%	81.3%
Sevenths	70.9%	67.4%	75.4%	79.7%	78.6%	84.1%
Intervals	67.0%	65.8%	69.7%	79.0%	77.9%	82.4%
Tonal Features from [45]						
	Composer Filter			No Filter		
	Full	*Piano*	*Orchestra*	*Full*	*Piano*	*Orchestra*
Template	73.9%	74.0%	79.4%	82.2%	84.9%	85.4%
Complexity	64.4%	68.1%	72.4%	76.6%	83.0%	82.1%
Complexity & Template	74.7%	73.3%	80.1%	85.0%	86.7%	87.5%

Table B.1: Further results of the style classification experiments.

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