

Similarity-Based Queries for Time Series Data

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Abstract

We study a set of linear transformations on the Fourier series representation of a sequence that can be used as the basis for similarity queries on time-series data. We show that our set of transformations is rich enough to formulate operations such as moving average and time warping. We present a query processing algorithm that uses the underlying R-tree index of a multidimensional data set to answer similarity queries efficiently. Our experiments show that the performance of this algorithm is competitive to that of processing ordinary (exact match) queries using the index, and much faster than sequential scanning. We relate our transformations to the general framework for similarity queries of Jagadish et al.

1 Introduction

Time-series data are of growing importance in many new database applications, such as data mining or warehousing. A time series is a sequence of real numbers, each number representing a value at a time point. For example, the sequence could represent stock or commodity prices, sales, exchange rates, weather data, biomedical measurements, etc. We are often interested in similarity queries on time-series data [APWZ95, ALSS95]. For example, we may want to find stocks that behave in approximately the same way (or approximately the opposite way, for hedging); or stocks that increased linearly up to October 1987, and then crashed; or years when the temperature patterns in two regions of the world were similar. In this type of queries, approximate rather than exact matching is required.

A naive approach is to compute the Euclidean distance (or any other distance, such as the city-block distance) between two objects (in general) or two time sequences (in particular), and call two sequences similar if their distance is less than a user-defined thresh-

old. Time sequences are usually long, so the distance computation can be time consuming. A solution is to map time sequences into the frequency domain using the Fourier transform, and use the first few coefficients to filter out non-similar data. This has the advantage that spatial indexing techniques can be used to index time sequences by viewing them as tuples of Fourier series coefficients, that is, points in a low-dimensional space [AFS93, FRM94].

A problem with this approach is that the user has no control over the meaning of similarity other than providing a threshold. There are many similarity queries that such a fixed predefined notion of similarity fails to capture; for example, one may consider two stocks similar if they have almost the same price fluctuations, even though one stock might sell twice as much as the other. Consider the following motivating examples:

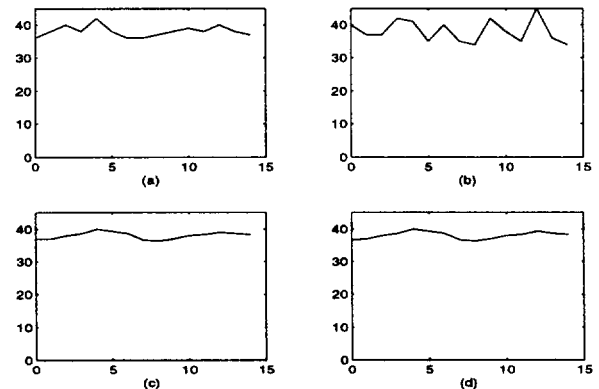


Figure 1: (a) Time sequence $s_1 = (36, 38, 40, 38, 42, 38, 36, 36, 37, 38, 39, 38, 40, 38, 37)$, (b) time sequence $s_2 = (40, 37, 37, 42, 41, 35, 40, 35, 34, 42, 38, 35, 45, 36, 34)$, (c) the 3-day moving average of s_1 , and (d) the 3-day moving average of s_2

Example 1.1 Suppose $s_1 = (36, 38, 40, 38, 42, 38, 36, 36, 37, 38, 39, 38, 40, 38, 37)$ and $s_2 = (40, 37, 37, 42, 41, 35, 40, 35, 34, 42, 38, 35, 45, 36, 34)$ are two time sequences that correspond to the closing prices of two stocks. Looking at Figure 1(a),(b), the sequences do not appear very similar. This is justified by the high Euclidean distance

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$D(\vec{s}_1, \vec{s}_2) = 11.92$ between them. However, if we look at the three-day moving averages of the two sequences (Figure 1 (c),(d)), they do look quite similar. The Euclidean distance between the three-day moving averages of two sequences is 0.47.

Moving averages are widely used in stock data analysis (for example, see [EM69]). Their primary use is to smooth out short term fluctuations and depict the underlying trend of a stock. The computation is simple; the l -day moving average of a sequence $\vec{s} = (v_1, \dots, v_n)$ is computed as follows: the mean is computed for an l -day-wide window placed over the end of the sequence; this will give the moving average for day $n - \lfloor l/2 \rfloor$; the subsequent values are obtained by stepping the window through the beginning of the sequence, one day at a time. This will produce a moving average of length $n - l + 1$. We use a slightly different version of moving average which is easier to compute in our framework. We circulate the window to the end of the sequence when it reaches the beginning. This gives us a moving average of length n . It turns out when the length of the window is small enough compared to the length of the sequence, which is usually the case in practice, both averages are almost the same.

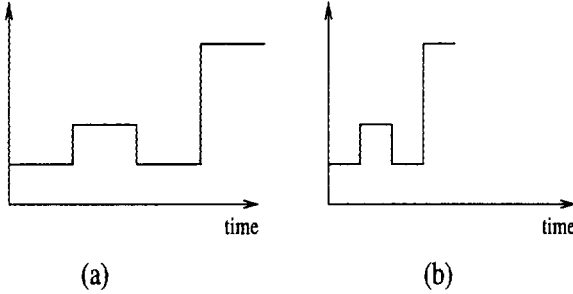


Figure 2: (a) Time sequence $\vec{s} = (20, 20, 21, 21, 20, 20, 23, 23)$ (b) time sequence $\vec{p} = (20, 21, 20, 23)$

Example 1.2 Consider two time sequences $\vec{s} = (20, 21, 21, 20, 21, 23, 23)$ and $\vec{p} = (20, 21, 20, 23)$ that are sampled with different frequencies (Figure 2). For example, \vec{s} could be the closing price of a stock taken every day, and \vec{p} could be the closing price of another stock taken every other day. A typical query is “is \vec{p} similar to \vec{s} ?”. The sequence \vec{s} is twice as long as \vec{p} , so they cannot be compared directly. The Euclidean distance between \vec{p} and any subsequence of length four of \vec{s} is more than 1.41. If the time dimension of \vec{p} is scaled by 2, i.e., every value “ v_i ” is replaced by “ v_i, v_i ”, the resulting sequence will be identical to \vec{s} . This operation is usually called *time warping* (for example, see [SK83]).

We propose a class of transformations that can be used in a query language to express similarity in a fairly general way, handling cases like the two examples above. Given an R-tree index [Gut84] constructed on a data set, we describe a fast query processing algorithm that uses the index to filter out unrelated data from the answer set of a similarity query. For example, we demonstrate

that an index structure for moving average can be constructed on the fly from the existing index, and it can be used to speed up the query processing. We show that this approach not only is faster than sequential scanning, but also introduces no extra disk overhead. To the best of our knowledge, this is the first indexing method that can handle moving average and time warping in the context of similarity queries.

The organization of the rest of the paper is as follows: The rest of the current section provides some basic material on the discrete Fourier transform and a survey of related work. Section 2 motivates the work by discussing possible applications to stock data analysis. Our definition of similarity queries is discussed in Section 3. In section 4 we develop an indexing method for these similarity queries. Section 5 presents experimental performance results. We conclude in Section 6.

1.1 The Discrete Fourier Transform

In this section, we briefly review the Discrete Fourier Transform (DFT) and its properties. Let a time sequence be a finite duration signal $\vec{x} = [x_t]$ for $t = 0, 1, \dots, n-1$. The DFT of \vec{x} , denoted by \vec{X} , is given by

$$X_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{n-1} x_t e^{-\frac{j2\pi ft}{n}} \quad f = 0, 1, \dots, n-1 \quad (1)$$

where $j = \sqrt{-1}$ is the imaginary unit. Throughout this paper, unless it is stated otherwise, we use small letters for sequences in the time domain and capital letters for sequences in the frequency domain. The inverse Fourier transform of \vec{X} gives the original signal, i.e.,

$$x_t = \frac{1}{\sqrt{n}} \sum_{f=0}^{n-1} X_f e^{\frac{j2\pi ft}{n}} \quad t = 0, 1, \dots, n-1 \quad (2)$$

Following the convention of [AFS93, FRM94], we have $1/\sqrt{n}$ in the front of both Equations 1 and 2. The energy of signal \vec{x} is given by the expression

$$E(\vec{x}) = \sum_{t=0}^{n-1} |x_t|^2. \quad (3)$$

The convolution of two signals \vec{x} and \vec{y} is given by

$$\text{Conv}(\vec{x}, \vec{y})_i = \sum_{k=0}^{n-1} x_k y_{i-k} \quad i = 0, 1, \dots, n-1 \quad (4)$$

where $i-k$ is computed modulo n . This convolution is usually called *circular convolution*. Equations 3 and 4 are unchanged in the frequency domain.

The following properties of DFT can be found in any signal processing textbook (for example, see [OS75]). The symbol \Leftrightarrow denotes a DFT pair.

Linearity

$$a\vec{x} + b\vec{y} \Leftrightarrow a\vec{X} + b\vec{Y} \quad (5)$$

for arbitrary constants a and b ,

Convolution-Multiplication

$$\text{conv}(\vec{x}, \vec{y}) \Leftrightarrow \vec{X} * \vec{Y} \quad (6)$$

where $\vec{X} * \vec{Y}$ is the element-to-element vector multiplication of two vectors \vec{X} and \vec{Y} , and

Parseval's Relation

$$E(\vec{x}) = E(\vec{X}). \quad (7)$$

Using Parseval's relation, it is easy to show that the Euclidean distance between two signals in the time domain is the same as their distance in the frequency domain.

$$D(\vec{x}, \vec{y}) = (E(\vec{x} - \vec{y}))^{1/2} = (E(\vec{X} - \vec{Y}))^{1/2} = D(\vec{X}, \vec{Y}) \quad (8)$$

A nice property of the DFT is that for a large family of sequences it concentrates the energy in the first few coefficients. Thus using the first few coefficients for indexing introduces few false hits, and no false dismissals.

1.2 Related Work

There has been some work on access methods for similarity queries. For example, Agrawal et al. [AFS93] propose an efficient index structure to retrieve time sequences similar to a given one. They map time sequences into the frequency domain using the Fourier transform and keep the first few coefficients in the index. Two sequences are considered similar if their Euclidean distance is less than a user-defined threshold. One difficulty with this approach is that the user has no control over the meaning of similarity.

Jagadish et al. [JMM95] develop a domain-independent framework to pose similarity queries on a database. The framework has three components: a *pattern language* P , a *transformation rule language* T , and a *query language* L . An expression in P specifies a set of data objects. An object A is considered similar to an object B , if B can be reduced to it by a sequence of transformations defined in T . The query language proposed in the paper is an extension of relational calculus with predicates that test whether an object A can be transformed into a member of the set of objects described by expression e using the transformation t , at a cost bounded by c . A specialization of this work to real-valued sequences where the search is performed over sequence signatures instead of the original sequences is described in [FJMM95].

In this paper, we describe an efficient implementation of a special case of [JMM95] for time-series data. We only study the trivial pattern language where a pattern expression specifies either a given constant data object, or every object in the database. Given an object o , a pattern expression e that denotes a set of objects, and a transformation t , the expression $t(e)$ ¹ denotes the set of all objects that can be obtained by applying t to every member of the set defined by e . We consider three kinds of queries: range queries, all-pair queries, and nearest neighbor queries, and we allow our transformations to be used in those queries.

We show how to use the indexing method in [AFS93] to test for similarity under a general class of transformations.

¹This is called $e \approx t$ in [JMM95].

Our work generalizes Goldin et al. [GK95] where transformations are limited to *shifts* and positive *scales*. Our extension allows shifts and scales in every dimension of a multidimensional feature space, as well as more complex transformations such as moving average. In addition, we drop the restriction to positive scales. Some advantages of these extensions are shown within the next section.

There are other related works on time series data. Agrawal et al. [APWZ95] describe a pattern language called SDL to encode queries about "shapes" found in time series. The language allows a kind of blurry matching where the user specifies the overall shape instead of the specific details, but it does not support any operations or transformations on the time series. A query language for time series data in the stock market domain is described in [Rot93]. The language is built on top of CORAL [RS92], and every query is translated into a sequence of CORAL rules.

2 Examples from Stock Data Analysis

In this section we demonstrate how our transformations can be used to eliminate noise or short-term fluctuations and shift or scale the data before computing Euclidean distances. We use three examples from real stock data. The data was obtained from the FTP site "ftp.ai.mit.edu/pub/stocks/results/".

Example 2.1 Figure 3 shows the daily closing price for *The Bombay Co.* (BBA) starting from October 25th, 1994 for 128 days, and that for *Zweig Total Return Fund Inc.* (ZTR) starting from July 20th, 1995 for 128 days. The Euclidean distance between two series is 16.16. The mean for BBA is 9.51, and the mean for ZTR is 8.64. If we shift the mean of both series to zero, i.e. subtract the mean of each series from everyday closing price, the Euclidean distance reduces to 12.78. The closing price of ZTR fluctuates in a smaller range than that of BBA; the standard deviation for ZTR is 0.10 while the standard deviation for BBA is 1.18. We scale both series by the inverse of their standard deviation. The resulting series in [GK95] are called *normal forms* of the original series. Thus given any sequence \vec{s} , sequence \vec{s}' is the normal form of \vec{s} if

$$s'_i = \frac{s_i - \text{mean}(\vec{s})}{\text{std}(\vec{s})} \quad \text{for } i = 1, \dots, \text{length}(\vec{s}) \quad (9)$$

Figure 3 shows that the Euclidean distance between the normal forms of two series is still 11.10; time series ZTR is more volatile than BBA. To smooth out short term fluctuations, we take the 20-day moving average of the two series. The Euclidean distance drops to 2.75.

Example 2.2 This example shows how we can identify series that have opposite price movements. Figure 4 shows the daily closing price for *Circuit City Stores Inc.* (CC) (marked by dotted lines) and the daily closing price for *Varian Associates Inc.* (VAR) (marked by solid lines) both starting from August 30, 1993 for 128 days. As Figure 4 shows, the two series have a reverse movement; when the price for CC goes up, the price for VAR goes down and vice versa. The Euclidean distance

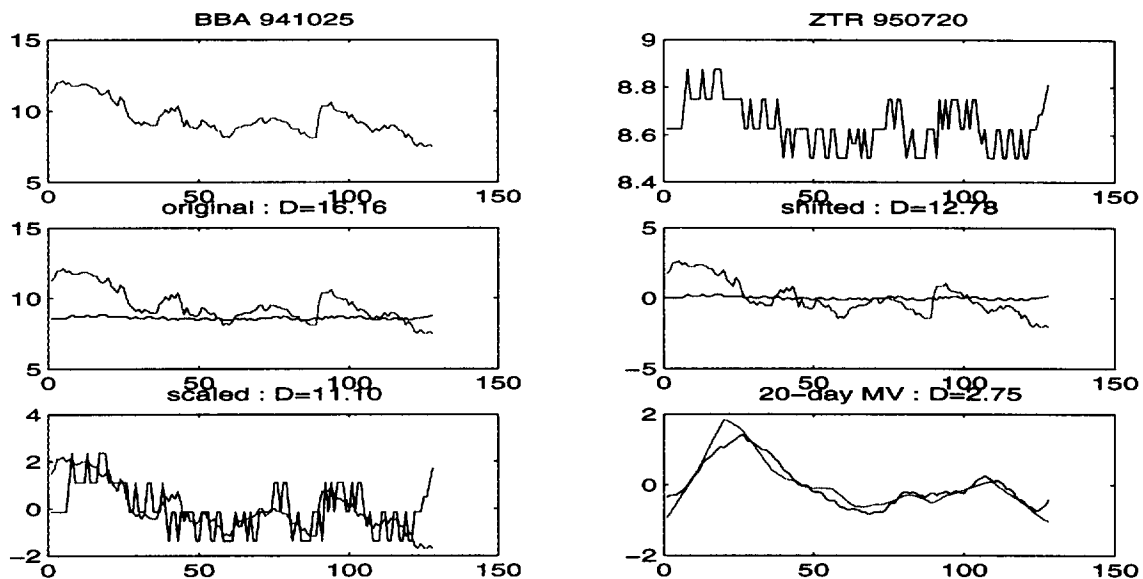


Figure 3: From left to right, top to bottom: the daily closing price for *The Bombay Co. (BBA)* starting from "94/10/25" for 128 days, the daily closing price for *Zweig Total Return Fund Inc. (ZTR)* starting from "95/07/20" for 128 days, the two stocks put together, both shifted, both scaled, and the 20-day moving average (D denotes the Euclidean distance)

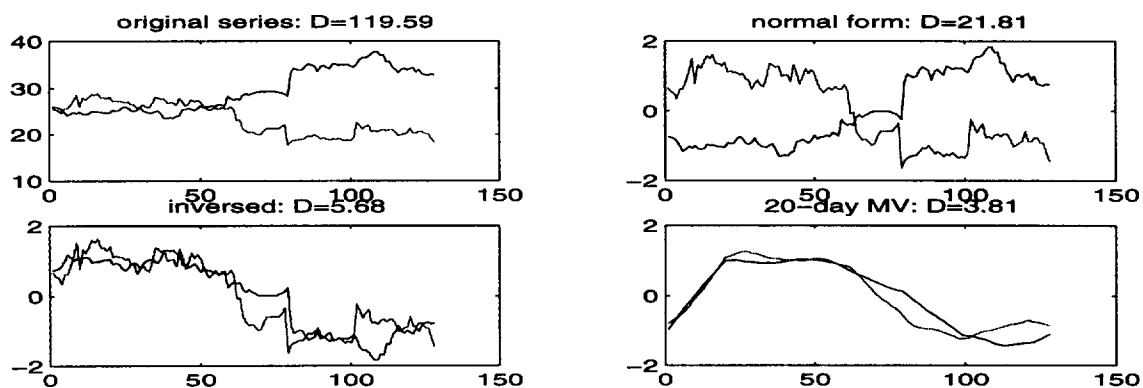


Figure 4: From left to right, top to bottom: the daily closing prices for *Circuit City Stores Inc. (CC)* (marked by dotted lines) and *Varian Associates Inc. (VAR)* (marked by solid lines) both starting from "93/08/30" for 128 days, their normal forms, VAR reversed, and the 20-day moving averages of both series (D denotes the Euclidean distance)

between two series is 119.59. We transform both series to their normal form, and the Euclidean distance becomes 21.81. If we reverse the time series of VAR, i.e., multiply everyday closing price by -1, and then take the 20-day moving average of both series, the Euclidean distance will be 3.81.

One might think that applying these transformations, any two series can be made similar. The next example shows this is not the case. Given three operations: *shift*, *scale*, and *20-day moving average*, we can use *shift* and *scale* to transform two series to their normal forms. We can smooth the normal form series using *20-day moving average*. Each one of these operations may reduce the distance between two series, but two series that have dissimilar trends still look different. It is obvious that if we keep taking the moving average, two series eventually will be the same, i.e., two flat straight lines. However, we assume, following [JMM95], that each operation has a cost, and we are limited by an upper bound on the total cost. This upper bound, for example, could be proportional to the Euclidean distance between the two original series.

Example 2.3 Figure 5 shows the daily closing prices of *Digital Microwave Corp. (DMIC)* and *The Mexico Fund Inc. (MXF)* both starting from August 30, 1993 for 128 days. The Euclidean distance between the normal forms of two series is 11.06. The Euclidean distance after taking the 20-day moving average becomes 10.09. The Euclidean distances after taking the second and the third 20-day moving average are respectively 9.63 and 9.22. The Euclidean distance even after taking the 10th moving average is still 6.57.

In the next section we encode the transformations discussed here in a query language.

3 Similarity Queries

We consider an object to be a point in a multidimensional space (md-space). For non-point objects, we assume there is a mapping function that maps every object to a point in the md-space. Such a function is developed in many domains where multidimensional indexing has been used. For example, Fourier transform for time-series [AFS93], and minimum bounding rectangle for shapes [Jag91] are some instances of the mapping function.

A transformation in an n -dimensional space, denoted by (\vec{a}, \vec{b}) , is a pair of $n \times 1$ vectors where \vec{a} specifies a stretch and \vec{b} represents a translation (Figure 6). The transformation (\vec{a}, \vec{b}) applied to a point \vec{x} in some space, maps \vec{x} to $\vec{a} * \vec{x} + \vec{b}$ which is a point in the same space. Transformations may be associated with costs. Given a set of transformations t , and the cost of applying each transformation, a measure of distance (dissimilarity) between two objects can be defined as follows:

$$D(\vec{x}, \vec{y}) = \min \begin{cases} D_0(\vec{x}, \vec{y}) \\ \min_{T \in t} (\text{cost}(T) + D(T(\vec{x}), \vec{y})) \\ \min_{T \in t} (\text{cost}(T) + D(\vec{x}, T(\vec{y}))) \\ \min_{T_1, T_2 \in t} (\text{cost}(T_1) + \text{cost}(T_2) + D(T_1(\vec{x}), T_2(\vec{y}))) \end{cases} \quad (10)$$

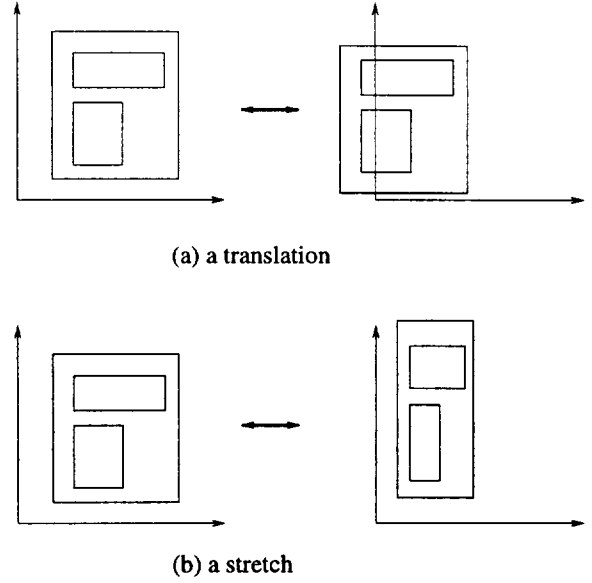


Figure 6:

where $D_0(\vec{x}, \vec{y})$ is the Euclidean distance between \vec{x} and \vec{y} . We use $T(\vec{x})$ to denote “transformation T applied to a point \vec{x} ” and $T(r)$ to denote “transformation T applied to a relation r ”. The former returns a point while the latter returns a relation. We assume relations are unary, that is, they are simply sets of sequences; in practice of course they may have other attributes, such as source of the data, time period covered, etc.

In the domain of time series data, both objects and transformations can be vectors of complex numbers, so we need to extend transformations to complex numbers. On the other hand, we want to make sure that this extension still keeps the main properties we are interested in. The following definition describes these properties.

Definition 1 A transformation T in a multidimensional space S is safe if T maps every rectangle R in S to a rectangle R' in S , every point inside R to a point inside R' , and every point outside R to a point outside R' .

Theorem 1 Transformation $T = (\vec{a}, \vec{b})$ is safe if \vec{a} and \vec{b} are chosen to be vectors of real numbers.

Proof (sketch): Transformation T here is the composition of a stretch and a translation in every dimension. Thus T is safe. ■

In the next section, we study the safety condition for complex numbers in more detail.

3.1 Transformations on Time Series

We consider a time series to be a point in a multidimensional feature space. We have chosen the first k Fourier coefficient of a time series as our features. The reason for choosing DFT is mainly because: (a) DFT concentrates the energy in the first few coefficients, so those coefficients can make a key for indexing purposes; (b) as we remarked in Section 1.1, it is known that the Euclidean distance is unchanged under the DFT. Since

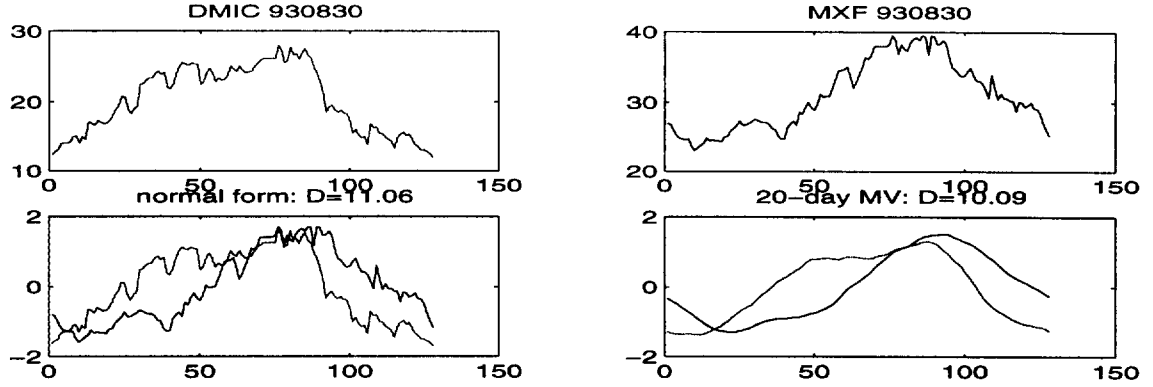


Figure 5: From left to right, top to bottom: the daily closing price of *Digital Microwave Corp.* (DMIC) starting from “93/08/30” for 128 days, the daily closing price of *The Mexico Fund Inc.* (MXF) for the same period, their normal forms, and the 20-day moving averages of both series (D denotes the Euclidean distance)

a Fourier coefficient, in general, is a complex number, we need a representation for complex numbers in our feature space. One possibility is to decompose a complex number into its real and imaginary components, and map each component to a dimension. For a given set of k features, we represent it with a point in a $2k$ -dimensional space. We denote the space built using this representation by S_{rect} . An alternative representation is to decompose a complex number into its components in the polar coordinate system. A complex number in the polar coordinate system is represented by a magnitude and a phase angle. We denote the space built using this representation by S_{pol} . We use $Re(x)$, $Im(x)$, $Abs(x)$, and $Angle(x)$ to denote respectively the real, the imaginary, the magnitude, and the phase angle of a complex number x . Now we need to show that the transformations described for time series data in previous sections are safe.

Theorem 2 Let \vec{a} be a vector of real numbers, and \vec{b} be a vector of complex numbers; the transformation $T = (\vec{a}, \vec{b})$ is safe with respect to S_{rect} .

Proof: We need to prove that T maps every rectangle R in the space to a rectangle R' , all interior points of R to interior points of R' , and all exterior points of R to exterior points of R' . Without loss of generality, we assume dimensions $2i-1$ and $2i$ (for $i = 1, \dots, k$) are respectively used for real and imaginary components of feature i . Suppose \vec{x} is a k -dimensional vector of complex numbers and \vec{x} , a $2k$ -dimensional vector, is its representation in S_{rect} . We have $x_i = x_{2i-1} + x_{2i}j$ for $i = 1, \dots, k$. If we apply transformation T to \vec{x} , we get $\vec{x}' = T(\vec{x}) = \vec{a} * \vec{x} + \vec{b}$. We can rewrite this as follows:

$$\begin{aligned} x'_i &= a_i * (x_{2i-1} + x_{2i}j) + (Re(b_i) + Im(b_i)j) \\ &= (a_i * x_{2i-1} + Re(b_i)) + (a_i * x_{2i} + Im(b_i))j \end{aligned}$$

for $i = 1, \dots, k$. If we map the resulting vector to a point x' in S_{rect} , we get $x'_{2i-1} = a_i * x_{2i-1} + Re(b_i)$ and $x'_{2i} = a_i * x_{2i} + Im(b_i)$ for $i = 1, \dots, k$. This transformation can be rewritten as $T' = (\vec{c}, \vec{d})$ where $c_{2i-1} = c_{2i} = a_i$, $d_{2i-1} = Re(b_i)$, and $d_{2i} = Im(b_i)$ for $i = 1, \dots, k$. Since

\vec{c} and \vec{d} are vectors of real numbers, the rest follows from Theorem 1. ■

On the other hand, Theorem 2 does not hold if \vec{a} is chosen to be a vector of complex numbers. For example consider a two dimensional rectangle with point $p = -5 - 5j$ as its lower left corner and point $q = 5 + 5j$ as its upper right corner, and $r = -2 + 2j$ as a point inside the rectangle. If we multiply the complex numbers representing the three points by $s = 2 - 3j$, the transformed rectangle built on points $p * s = -25 + 5j$ and $q * s = 25 - 5j$ does not have point $r * s = 2 + 10j$ inside!

Theorem 3 Let \vec{a} be a vector of complex numbers, and \vec{b} be a zero vector ($\vec{b} = \vec{0}$); the transformation $T = (\vec{a}, \vec{b})$ is safe with respect to S_{pol} .

Proof: Without loss of generality, we assume dimensions $2i-1$ and $2i$ (for $i = 1, \dots, k$) are respectively used for magnitude and phase angle of feature i . Suppose \vec{x} is a k -dimensional vector of complex numbers and \vec{x} is its coordinate in S_{pol} . We have $x_i = x_{2i-1}e^{x_{2i}j}$ for $i = 1, \dots, k$. If we apply transformation T to \vec{x} , we get $\vec{x}' = T(\vec{x}) = \vec{a} * \vec{x} + \vec{b}$. We can rewrite this as follows:

$$\begin{aligned} x'_i &= Abs(a_i)e^{Angle(a_i)j} * x_{2i-1}e^{x_{2i}j} + 0 \\ &= (Abs(a_i) * x_{2i-1})e^{(x_{2i} + Angle(a_i))j} \end{aligned}$$

for $i = 1, \dots, k$. If we map the resulting vector to a point x' in S_{pol} , we get $x'_{2i-1} = Abs(a_i) * x_{2i-1}$ and $x'_{2i} = x_{2i} + Angle(a_i)$ for $i = 1, \dots, k$. This transformation can be rewritten as $T' = (\vec{c}, \vec{d})$ where $c_{2i-1} = Abs(a_i)$, $d_{2i-1} = 0$, $c_{2i} = 1$, and $d_{2i} = Angle(a_i)$ for $i = 1, \dots, k$. Since \vec{c} and \vec{d} are vectors of real numbers, the rest follows from Theorem 1. ■

Given a query point \vec{q} in a $2k$ dimensional space and a threshold ϵ , we need to build a search rectangle. A search rectangle is the minimum bounding rectangle that contains all points within the Euclidean distance ϵ from \vec{q} . It is straightforward in the rectangular coordinate system; the minimum bounding rectangle is $(q_i - \epsilon, q_i + \epsilon)$ for $i = 1, \dots, 2k$. The minimum bounding

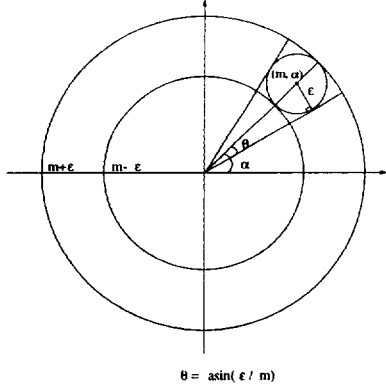


Figure 7: Minimum bounding rectangle in the polar coordinate system

rectangle for a complex number me^{aj} in the polar coordinate system is demonstrated in Figure 7. The magnitude is in the range from $m-\epsilon$ to $m+\epsilon$, and the angle is in the range from $\alpha - \text{asin}(\frac{\epsilon}{m})$ to $\alpha + \text{asin}(\frac{\epsilon}{m})$ where asin denotes the arc sinus of an angle. If we again assume dimensions $2i-1$ and $2i$ (for $i = 1, \dots, k$) are respectively used for magnitude and phase angle of feature i , then the minimum bounding rectangle for \vec{q} in the polar coordinate will be $(q_i - \epsilon, q_i + \epsilon)$ for $i = 1, 3, 5, \dots, 2k-1$ and $(q_i - \text{asin}(\frac{\epsilon}{q_{i-1}}), q_i + \text{asin}(\frac{\epsilon}{q_{i-1}}))$ for $i = 2, 4, 6, \dots, 2k$.

3.2 Using Transformations to Express Similarities

To gain some insight into the transformations, we formalize the notion of similarity expressed in Example 1.1. Time series \vec{s}_1 is considered similar to \vec{s}_2 because their 3-day moving averages look the same, so we need to formulate the 3-day moving average in our transformation language. For simplicity, in our examples we assign a cost of zero to all transformations. Let us denote the Fourier transform of \vec{s}_1 by \vec{S}_1 , the Fourier transform of \vec{s}_2 by \vec{S}_2 , and the Fourier transform of $\vec{m}_3 = (\frac{1}{3}, \frac{1}{3}, \frac{1}{3}, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0)$ by \vec{M}_3 . Now consider the transformation $T_{avg3} = (\vec{M}_3, \vec{0})$ where $\vec{0}$ is a zero vector of the same size as \vec{M}_3 . If we apply the transformation T_{avg3} to \vec{S}_1 , i.e.,

$$T_{avg3}(\vec{S}_1) = \vec{S}_1 * \vec{M}_3 + \vec{0} = \vec{S}_1 * \vec{M}_3$$

we get the 3-day moving average of \vec{s}_1 in the frequency domain. If we transform the right hand side back to the time domain using the convolution-multiplication relation (Equation 6), we get $T_{avg3}(\vec{s}_1) = \text{conv}(\vec{s}_1, \vec{m}_3)$ which is the 3-day moving average of \vec{s}_1 in the time domain. The same transformation can be applied to \vec{s}_2 to compute its 3-day moving average.

The notion of similarity can be expressed in a query by the proper choice of transformations. For example, the m -day moving average of a series of length n gener-

ally can be expressed by $T_{avg} = (\vec{a}, \vec{0})$ where

$$\vec{a} = (\underbrace{w_1 = 1/m, w_2 = 1/m, \dots, w_m = 1/m}_{m}, \underbrace{0, 0, \dots, 0}_n) \quad (11)$$

and $\vec{0}$ is a zero vector of size n . Transformation T_{avg} may be applied several times to get successive moving averages. The weights w_1, \dots, w_m are not necessarily equal. For trend prediction purposes, for example, the weights at the end are usually chosen to be higher than those at the beginning. Whereas for normal smoothing purposes, weights are equal, or those at the center are larger than those at the endpoints.

To give another example of the transformations, we formulate the transformation used to reverse a time series in Example 2.2. Let $T_{rev} = (\vec{a}, \vec{0})$ where $a_i = -1$ for $i = 1, \dots, 128$ and $\vec{0}$ is a zero vector of size 128. Now consider a time series \vec{s} and its Fourier transform \vec{S} . Transformation T_{rev} applied to \vec{S} gives

$$T_{rev}(\vec{S}) = \vec{a} * \vec{S} + \vec{0} = -\vec{S}.$$

If we transform both sides into the time domain using Equation 5, we get $T_{rev}(\vec{s}) = -\vec{s}$. That is, the daily closing price is multiplied by -1 .

Transformation T_{rev} can be used to obtain all the pairs of series that move in opposite directions. This can be formulated in our query language for a given relation r as a spatial join between r and $T_{rev}(r)$.

Transformations can also be defined to stretch the time dimension (Example 1.2). Details of this transformation are given in Appendix A. In the next section, we discuss an indexing technique for similarity queries.

4 Indexing of Similarity Queries

In this section we describe a fast query processing method for similarity queries. We assume a multidimensional index is available, and we take an advantage of that in our query processing. Because of the dominant use of the R-tree family in multidimensional indexing, we describe our approach for R-tree indexes.

Given an R-tree index I in a multidimensional space S over a data set D , and any safe transformation T in S , we give an algorithm to construct an R-tree index I' for $T(D)$.

Algorithm 1 : For every node n

$$n = ((MBR_1, \text{pointer}_1), \dots, (MBR_m, \text{pointer}_m))$$

in I , we construct a node n'

$$\begin{aligned} n' &= T(n) \\ &= ((MBR'_1, \text{pointer}'_1), \dots, (MBR'_m, \text{pointer}'_m)) \end{aligned}$$

in I' such that $MBR'_i = T(MBR_i)$, and $\text{pointer}'_i$ is a pointer to $T(n_i)$ where n_i is the child node (or the data tuple when n is a leaf node) pointed by pointer_i (for $i = 1, \dots, m$). We assumed T is a safe transformation, therefore MBR'_i is a bounding rectangle for all rectangles of the child node (or the data tuple) $T(n_i)$. The construction stops when every node in I is mapped to a node in I' .

There are many possible R-tree indexes on $T(D)$, each with a different performance. Our experiments show that the index we build here has a similar performance to that of the original index. The main observation here is that for a given index I and transformation T , index I' can be built on the fly without having much impact on the performance of the search. This allows us to use one index for many transformations.

An indexing method for time series data is described in [AFS93]. This method requires a cut-off point for the number of Fourier coefficients kept in the index. We denote this cut-off point by k and call the index built on the first k Fourier coefficients ' k -index'.

To demonstrate the query processing algorithm, we use a more general form of Example 1.1 throughout this section. We have seen in Section 3 that the m -day moving average of a series can be expressed by $T_{avg} = (\vec{a}, \vec{0})$ where \vec{a} is a vector of complex numbers. Due to Theorem 3, T_{avg} is safe if we represent complex numbers in the polar coordinate.

Query: Given a pattern expression e , a safe transformation T , an object \vec{q} , and a threshold ϵ , find all objects $\vec{o} \in T(e)$ such that the Euclidean distance $D(\vec{o}, \vec{q}) < \epsilon$.

If the pattern expression e denotes only one object \vec{o}_1 , we simply apply T to \vec{o}_1 . Object \vec{o}_1 is in the answer set if $D(\vec{o}_1, \vec{q}) < \epsilon$. Now suppose the expression e denotes all objects in the relation. A naive evaluation requires reading the whole relation, applying T to every object, and choosing every object \vec{o} such that $D(\vec{o}, \vec{q}) < \epsilon$. This is a costly process. A better approach is to use Algorithm 1 to construct a new index for transformed objects, and this new index can be built on the fly during the search operation. The search algorithm for the given range query is as follows:

Algorithm 2 : Given a k -index whose root is N , a transformation T , a threshold ϵ , and a search point \vec{q} , apply T to all points in the index and find those whose distance from \vec{q} becomes less than ϵ .

1. Preprocessing:

- (a) Transform T and \vec{q} into the frequency domain if they are in the time domain. Let us denote the first k Fourier coefficients of T and \vec{q} by T_k and \vec{q}_k respectively.
- (b) build a search rectangle q_{rect} for \vec{q}_k as described in Section 3.1.

2. Search:

- (a) If N is not a leaf, apply T to every (rectangle) entry of N and check if the resulting rectangle overlaps q_{rect} . For all overlapping entries, call **Search** on the index whose root node is pointed to by the overlapping entry.
- (b) If N is a leaf, apply T to every (point) entry of N and check if the resulting point overlaps q_{rect} . If so, the entry is a candidate.

3. Postprocessing:

- (a) For every candidate point, check its full database record to determine if its Euclidean distance is at most ϵ from \vec{q} . If so, the entry is in the answer set.

Similarly all-pairs queries and nearest neighbor queries can be processed efficiently using the index. For an all-pairs query, we do a spatial join using the index. The only difference here is that we transform all objects used in the join predicate before we compute the predicate. For example, the join predicate $a_i \cap b_j \neq \emptyset$ may be changed to $T(a_i) \cap T(b_j) \neq \emptyset$ where T is a transformation and a_i and b_j are members of two spatial sets. For a nearest neighbor query, the search starts from the root and proceeds down the tree. As we go down the tree, we apply T to all entries of the node we visit. We can then use any kind of metric (such as MINDIST or MINMAXDIST discussed in [RKV95]) for pruning the search.

The only thing left to show is that this search scheme used with a k -index misses no object from the answer set.

Lemma 1 *The k -index approach enhanced with transformations always returns a superset of the answer set.*

Proof: Suppose we want to find all objects \vec{x} in a relation that are similar to a query object \vec{q} . Since transformations are applied to series in the frequency domain, this can be written in the frequency domain as follows:

$$D(T(\vec{X}), \vec{Q}) \leq \epsilon \quad (12)$$

where $T = (\vec{A}, \vec{B})$ is a transformation, ϵ is a user-defined threshold, and \vec{X} and \vec{Q} are DFTs of respectively \vec{x} and \vec{q} . Applying T to \vec{X} , we get

$$D(\vec{A} * \vec{X} + \vec{B}, \vec{Q}) = \left(\sum_{f=0}^{n-1} |A_f X_f + B_f - Q_f|^2 \right)^{\frac{1}{2}} \leq \epsilon$$

If we keep only the first $k < n$ coefficients, we have

$$\left(\sum_{f=0}^{k-1} |A_f X_f + B_f - Q_f|^2 \right)^{\frac{1}{2}} \leq \left(\sum_{f=0}^{n-1} |A_f X_f + B_f - Q_f|^2 \right)^{\frac{1}{2}} \quad (13)$$

On the other hand, the equation

$$\left(\sum_{f=0}^{n-1} |A_f X_f + B_f - Q_f|^2 \right)^{\frac{1}{2}} \leq \epsilon \quad (14)$$

holds for all objects in the answer set. Equations (13,14) imply that

$$\left(\sum_{f=0}^{k-1} |A_f X_f + B_f - Q_f|^2 \right)^{\frac{1}{2}} \leq \epsilon \quad (15)$$

Therefore, keeping the first k coefficients introduces no false dismissals. ■

This is a generalization of the result of [AFS93] for k -index enhanced with transformations.

5 Experiments

We implemented our method on top of Norbert Beckmann's Version 2 implementation of the R*-tree [BKSS90]. We ran experiments on both stock prices data obtained from the FTP site "ftp.ai.mit.edu/pub/stocks/results/" and synthetic sequences. Each synthetic sequence $\vec{x} = [x_i]$ was a random sequence produced as follows:

$$\begin{aligned} x_0 &= y \\ x_1 &= x_0 + z_1 \\ &\dots \\ x_i &= x_{i-1} + z_i \end{aligned}$$

where y was a normally distributed random number in the range $[20, 99]$, and z_i ($i = 1, 2, \dots$) was a random number in the range $[-4, 4]$.

We used the polar representation of complex numbers because vector multiplication for time series data seemed to be more important than vector addition, and due to Theorem 3 vector multiplication is safe with respect to S_{pol} . For every time series, we first transformed it to the normal form, and then we found its Fourier coefficients. The reason for choosing the normal form was because both the average and the standard deviation of a series could be stored in the index as two separate dimensions, and despite using the polar representation, we could still have simple shifts. Since the mean of a normal form series is zero by definition, the first Fourier coefficient is always zero, so we can throw it away. We mapped the mean and the standard deviation of the original time series respectively to the first and the second dimensions of the index. We also mapped the magnitude and the phase angle of the second DFT term (computed for the normal form series) respectively to the third and the fourth dimensions of the index, and the magnitude and the phase angle of the third DFT term respectively to the fifth and the sixth dimensions.

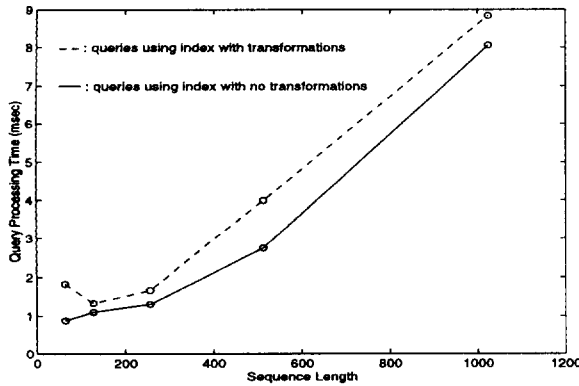


Figure 8: time per query varying the sequence length

Figure 8 compares the execution time for two kinds of queries: (a) a range query using transformations and (b) a range query that uses no transformations. We varied the length of the sequences from 64 to 1024 while we kept the number of sequences fixed to 1,000. In order to have a precise comparison, the identity transformation

$T_i = (\vec{I}, \vec{0})$ was chosen such that $T_i(\vec{o}) = \vec{o}$ for all objects \vec{o} (\vec{I} is a vector of 1's and $\vec{0}$ is a vector of 0's). This made the two queries produce the same results. As Figure 8 shows the difference between the two curves is only a constant. This constant is the CPU time spent for vector multiplication which is unavoidable. The number of disk accesses is the same in both cases.

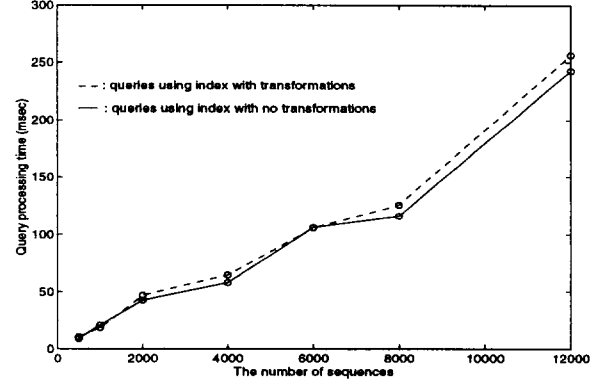


Figure 9: time per query varying the number of sequences

In the next experiment, we kept the sequence length fixed to 128 while we varied the number of sequences from 500 to 12,000. We used the identity transformation again for the same reason described in the previous experiment. As demonstrated in Figure 9, the result was the same. Thus the index traversal for similarity queries does not deteriorate the performance of the index.

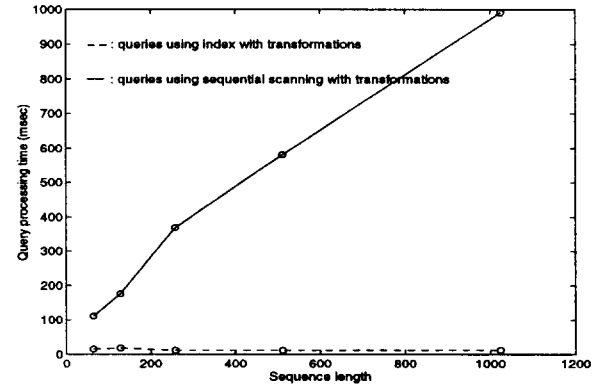


Figure 10: time per query varying the sequence length

Figures 10 and 11 compare the execution time of our approach to sequential scanning. To have a good implementation for the sequential scan, we stop the distance computation process as soon as the distance exceeds ϵ . In addition, we do the sequential scanning on the relation that stores the series in the frequency domain, not the time domain. Because each series in the frequency domain has its larger coefficients at the beginning, the distance computation process can skip many sequences within the first few coefficients. Both graphs show the superiority of our approach.

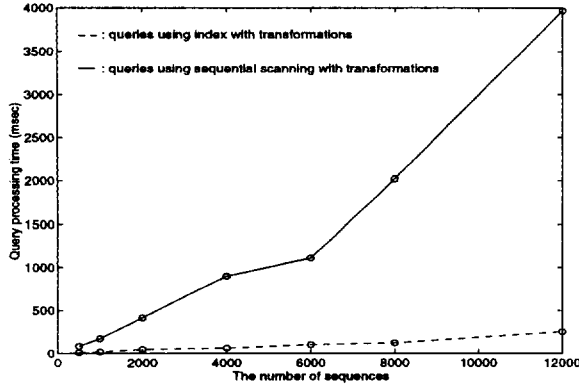


Figure 11: time per query varying the number of sequences

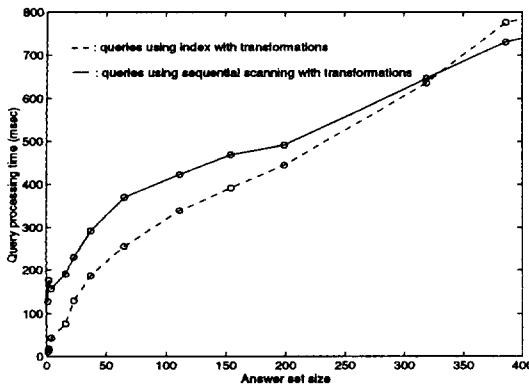


Figure 12: time per query varying the size of the answer set

In another experiment, we kept the number of sequences fixed to 1067, and we also kept the sequence length fixed to 128. The experiment ran on the real stock data. We varied the threshold so that the query gave us different numbers of time series in the answer set. Figure 12 shows that the index performs better until the size of the answer set gets larger than 300 which is almost one third of the size of the relation.

Our last experiment was on a spatial self-join. In fact, all the time series used as examples in Section 2 were the results of this spatial join. We did the test using the following methods:

- a scan the relation of Fourier coefficients sequentially, and compare every sequence s to all the sequences that are after s in the relation; the transformation T_{avg20} is applied to every sequence during the comparison;
- b do the sequential scanning as instructed in a, but stop the distance computation as soon as the distance exceeds ϵ .
- c scan the relation of Fourier coefficients sequentially, and for every sequence build a search rectangle and pose it to the index as a range query;
- d do the spatial join as described in c, but in every index retrieval apply T_{avg20} to both the index and the search rectangles.

The experiment ran on a relation of stock prices data that had 1067 time sequences, and the length of each sequence was 128. The result of the test is shown in table 1. Because of the implementation, b is 10 times

algorithm	time (min:sec.milise)	size of the answer set
a	20:36.323	12
b	2:31.217	12
c	0:10.139	$3 \times 2 = 6$
d	0:17.698	$12 \times 2 = 24$

Table 1: The result of the join

faster than a. Methods c and d are 9 to 15 times faster than b because of using the index, and d is a bit slower than c because of the use of a transformation and the size of the answer set. The answer set of d contains every pair twice, so it is twice the size of a and b. The size of the answer set for c is smaller because it does not use the transformation.

6 Conclusions

We have proposed a class of transformations that can be used in a query language to express similarity among objects. This class allows the expression of several practically important notions of similarity, and queries using this class can be efficiently implemented on top of any R-tree index. One potential application which is emphasized in the paper is stock data analysis, but we believe other domains can also benefit. Our contributions can be summarized as follows:

- Formulation of moving average, time warping, and reversing in our transformation language.
- Implementation of similarity matching under these transformations on top of an R-tree index.

The experiments show that execution time of our method is almost the same as that of accessing the index with no transformations; our method has much better performance than sequential scanning, and the performance gets better by increasing both the number and the length of sequences.

We think the normal form of [GK95] is a useful representation for time series data, but it allows only a small fraction of similarity queries. Our similarity transformations allow more general queries, but for simple shifting and scaling, the indexing method in [GK95] is faster because no transformation needs to be performed on the index. Our indexing technique can be easily built on top of [GK95] as we did in our experiments, allowing both simple shifts and scales and more general transformations to be applied efficiently.

Acknowledgments

We would like to thank Christos Faloutsos for his help in providing us with stock data and his comments on a preliminary version of this paper. This work was supported by the Natural Sciences and Engineering Research Council of Canada and the Information Technology Research Centre of Ontario.

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A Time Warping

Given the first $k \leq n$ Fourier coefficients of a time series \vec{s} of length n and an integer $m \geq 1$, we can construct the first k Fourier coefficients of time series \vec{s}' of length $m \times n$ using a transformation $T = (\vec{a}, \vec{0})$ such that

$$s'_{mi} = s'_{mi+1} = \dots = s'_{m(i+1)-1} = s_i \quad (16)$$

for $i = 0, \dots, n$.

Let $\vec{s} = (s_1, s_2, \dots, s_n)$ be a time sequence, and $\vec{S} = (S_1, S_2, \dots, S_n)$ be its DFT. Using Equation 1, we can write

$$S_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{n-1} s_t e^{\frac{-j2\pi tf}{n}} \quad f = 0, 1, \dots, n-1. \quad (17)$$

We want to find a vector \vec{a} such that

$$a_f * S_f = S'_f \quad \text{for } f = 0, \dots, k-1. \quad (18)$$

where S'_f is the f th Fourier coefficient of \vec{S}' . Using Equation 1, we can write S'_f as follows:

$$S'_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{mn-1} s'_t e^{\frac{-j2\pi tf}{mn}}.$$

This can be rewritten as

$$\begin{aligned} S'_f = \frac{1}{\sqrt{n}} & \left(\sum_{t=0}^{m-1} s'_t e^{\frac{-j2\pi tf}{mn}} + \sum_{t=m}^{2m-1} s'_t e^{\frac{-j2\pi tf}{mn}} + \dots \right. \\ & \left. + \sum_{t=(n-1)m}^{nm-1} s'_t e^{\frac{-j2\pi tf}{mn}} \right). \end{aligned}$$

If we rewrite all summations as $\sum_{t=0}^{m-1}$ and also use Equation 16, we get

$$\begin{aligned} S'_f = \frac{1}{\sqrt{n}} & \left(\sum_{t=0}^{m-1} s_0 e^{\frac{-j2\pi tf}{mn}} + \sum_{t=0}^{m-1} s_1 e^{\frac{-j2\pi(t+m)f}{mn}} + \dots \right. \\ & \left. + \sum_{t=0}^{m-1} s_{n-1} e^{\frac{-j2\pi(t+(n-1)m)f}{mn}} \right). \end{aligned}$$

We can take $\sum_{t=0}^{m-1} e^{\frac{-j2\pi tf}{mn}}$ out of the parenthesis; that gives

$$S'_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{m-1} e^{\frac{-j2\pi tf}{mn}} (s_0 + s_1 e^{\frac{-j2\pi f}{n}} + \dots + e^{\frac{-j2\pi(n-1)f}{n}}).$$

Using Equation 17, we can rewrite this equation as follows:

$$S'_f = \sum_{t=0}^{m-1} e^{\frac{-j2\pi tf}{mn}} S_f.$$

Therefore, if we choose vector \vec{a} as follows:

$$a_f = \sum_{t=0}^{m-1} e^{\frac{-j2\pi tf}{mn}} \quad \text{for } f = 0, \dots, k-1, \quad (19)$$

then the Equation 18 holds.

A Time Warping

Given the first $k \leq n$ Fourier coefficients of a time series \vec{s} of length n and an integer $m \geq 1$, we can construct the first k Fourier coefficients of time series \vec{s}' of length $m \times n$ using a transformation $T = (\vec{a}, \vec{0})$ such that

$$s'_{mi} = s'_{m(i+1)} = \dots = s'_{m(i+1)-1} = s_i \quad (16)$$

for $i = 0, \dots, n$.

Let $\vec{s} = (s_1, s_2, \dots, s_n)$ be a time sequence, and $\vec{S} = (S_1, S_2, \dots, S_n)$ be its DFT. Using Equation 1, we can write

$$S_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{n-1} s_t e^{-j2\pi t f / n} \quad f = 0, 1, \dots, n-1. \quad (17)$$

We want to find a vector \vec{a} such that

$$a_f * S_f = S'_f \quad \text{for } f = 0, \dots, k-1. \quad (18)$$

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$$S'_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{mn-1} s'_t e^{-j2\pi t f / mn}.$$

This can be rewritten as

$$S'_f = \frac{1}{\sqrt{n}} \left(\sum_{t=0}^{m-1} s'_t e^{-j2\pi t f / mn} + \sum_{t=m}^{2m-1} s'_t e^{-j2\pi t f / mn} + \dots \right. \\ \left. + \sum_{t=(n-1)m}^{nm-1} s'_t e^{-j2\pi t f / mn} \right).$$

If we rewrite all summations as $\sum_{t=0}^{m-1}$ and also use Equation 16, we get

$$S'_f = \frac{1}{\sqrt{n}} \left(\sum_{t=0}^{m-1} s_0 e^{-j2\pi t f / mn} + \sum_{t=0}^{m-1} s_1 e^{-j2\pi (t+m) f / mn} + \dots \right. \\ \left. + \sum_{t=0}^{m-1} s_{n-1} e^{-j2\pi (t+(n-1)m) f / mn} \right).$$

We can take $\sum_{t=0}^{m-1} e^{-j2\pi t f / mn}$ out of the parenthesis; that gives

$$S'_f = \frac{1}{\sqrt{n}} \sum_{t=0}^{m-1} e^{-j2\pi t f / mn} (s_0 + s_1 e^{-j2\pi f / n} + \dots + e^{-j2\pi (n-1) f / n}).$$

Using Equation 17, we can rewrite this equation as follows:

$$S'_f = \sum_{t=0}^{m-1} e^{-j2\pi t f / mn} S_f.$$

Therefore, if we choose vector \vec{a} as follows:

$$a_f = \sum_{t=0}^{m-1} e^{-j2\pi t f / mn} \quad \text{for } f = 0, \dots, k-1, \quad (19)$$

then the Equation 18 holds.