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Developing a Pattern Discovery Method in Time Series Data and its GPU Acceleration

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Abstract: The Dynamic Time Warping (DTW) algorithm is widely used in finding the global alignment of time series. Many time series data mining and analytical problems can be solved by the DTW algorithm. However, using the DTW algorithm to find similar subsequences is computationally expensive or unable to perform accurate analysis. Hence, in the literature, the parallelisation technique is used to speed up the DTW algorithm. However, due to the nature of DTW algorithm, parallelising this algorithm remains an open challenge. In this paper, we first propose a novel method that finds the similar local subsequence. Our algorithm first searches for the possible start positions of subsequence, and then finds the best-matching alignment from these positions. Moreover, we parallelise the proposed algorithm on GPUs using CUDA and further propose an optimisation technique to improve the performance of our parallelization implementation on GPU. We conducted the extensive experiments to evaluate the proposed method. Experimental results demonstrate that the proposed algorithm is able to discover time series subsequences efficiently and that the proposed GPU-based parallelization technique can further speedup the processing.

Key words: Dynamic Time Warping, Time series data, Data mining, Pattern Discovery, GPGPU, Parallel processing

1 Introduction

Time series data is a series of values sampled at time intervals [1]. Time series data captures a lot of information including inter-patterns, trends, and correlations. Also, time series data usually contains a high volume of numerical data and a time dimension. Many different techniques have been developed to

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model, compare, and predict time series data.

Due to the ubiquity of time series data, many real life data mining tasks can be modelled as mining time series data. These jobs often relate to finding known patterns, discovering unknown patterns and locating common subsequences in time series data [2–5]. Locating common subsequence or sequence matching is an important domain in time series data analysis.

As suggested in [6] and [7], many applications can be solved by finding the common subsequence pairs in two time series data. For example: finding the association rules in time series data [8,9]; classification algorithms that are based on building typical prototypes of each class [10, 11]; anomaly detection [12] and finding periodic patterns [13].

Dynamic Time Warping (DTW)[14] is a well-known algorithm for comparing similarities between two time series data. DTW algorithm is designed for finding the global sequence alignment in speech recognition.

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To find local alignments, previous researchers focus on finding subsequence within a longer sequence that match a shorter query sequence (Subsequence DTW) [15–17].

However, DTW algorithm suffers from high computational complexity. A straightforward implementation of DTW is quadratic in time and space. Using DTW to find local similarity between two time series normally requires applying the DTW algorithm to input data multiple times [18–20]. Therefore, it is necessary to improve the performance of the DTW algorithm.

Based on the above considerations, many efforts are made to develop methods and techniques that execute the DTW algorithm in parallel to reduce the execution time. Recently, with the development of general purpose GPUs and programming environment, much research is conducted to study how to execute the DTW algorithm on GPUs [21–23].

However, due to the data dependency, parallelising DTW algorithm on GPU is challeging and suffers from low parallelisation degree [21]. Therefore, in this paper, we propose a novel method to parallelise the DTW algorithm on GPU. Based on dynamic programming, we first propose a new algorithm of discovering similar local subsequences between time series sequences efficiently. Then we parallelise the proposed algorithm on GPU and use the diagonal order technique to optimise the memory access pattern in GPU.

This paper is organised as follows: Section 2 and 3 provide the related work and the backgrounds in time series data analysis. Section 4 proposes the design of the subsequence matching method. Section 5 presents the GPU implementation of the parallelised algorithm. We present the evaluation and experimental results in Section 6. Section 7 concludes this paper.

2 Related Work

Many researchers have conducted work in finding subsequences in time series data. In [6], Lin et al. developed a symbolic representation-based local subsequence matching algorithm to find the k^{th} most similar subsequence, but the algorithm requires a userinput length of the subsequence. A probabilistic method for subsequence matching is proposed in [7]. The method tries to match arbitrary subsequence of one time series to an arbitrary subsequence of another time series and analyze the statistical pattern of a result matrix to locate the position of the sequence. This method also requires the length of the subsequence to be known in advance. In [24], a local subsequence matching algorithm for multi-dimensional data is proposed. [25] presents a method of discovering exact subsequence (the length and the shape of the subsequence are known in advance). The algorithm provides a feasible solution in finding exact subsequence from short time series in a reasonable time. However, the time complexity of the method is still very high, which makes it not practical to work with long time series.

Toyoda et al. [26] proposed a novel method to discover similar subsequence in time series. The algorithm creates two matrices, score matrix and position matrix respectively. The score matrix locates the end position of the found subsequence, and the position matrix uses the information collected from the process of computing score matrix to calculate the start position of the subsequence. This method can find subsequences at any length and of any shape without producing meaningless subsequences and reduces required time and space complexity significantly in the meantime. Moreover, this method only needs three parameters and does not require prior knowledge of the subsequences to be found. However, this algorithm is still lack of scalability thanks to the exponential increase of matrix size with the accumulation of time series.

The work in [27] exploits the possibility of using CPU clusters to speed up DTW. They assign subsequences starting from different positions of the time series to different processors, and each processor performs the classic DTW algorithm. The method proposed in [28] explores the use of the multi-core processor to parallelize the DTW algorithm. In the work, they separate different patterns into different cores, and each subsequence will be assigned to different cores, where it is compared with different patterns naively. In these two parallel implementations, because a subsequence consists of several hundred or thousands of tuples, the data transferring becomes the bottleneck of the system.

In [23] a GPU implementation of DTW algorithm is proposed. In this work, multi-threads are used to generate the warping matrix in parallel, but the path search is performed serially. This implementation exploits the partial fine-grained parallelism of DTW itself. Because it separates the matrix generation phase and the path search phase into two kernels, it needs to store and transfer the whole warping matrix for each DTW calculation, which is also a heavy burden in the whole system. In [22] the authors claim that they are "the first to present hardware acceleration techniques for similarity search in streams under the DTW measure". Their GPU implementation is similar to [27].

3 Background

In this section we review the concepts of time series and dynamic time warping (DTW) distance [14] as well as some background knowledge for the ease of further discussion.

3.1 Time series

A time series can be represented as $T = (t_1, t_2, \ldots, t_n)$ which denotes ordered values in a time series of n samples [14]. The sequence are measured within fixed time intervals. A subsequence S with length k of T is denoted as $S = (t_i, t_{i+1}, \ldots, t_{i+k})$, where $1 \le i \le n - k$.

Given sequence $X = (x_1, x_2, \ldots, x_n)$ of length n and sequence $Y = (y_1, y_2, \ldots, y_m)$ of length m, the similar subsequences of X and Y are S_x and S_y that satisfy the pre-defined similarity measurement condition. In this paper, we aim to find all similar subsequences X_s and Y_s from two time series.

3.2 Classic DTW algorithm

Since the classic Dynamic Time Warping (DTW) algorithm serves as the foundation of our proposed algorithm in this paper, we introduce the algorithm in this section briefly.

The classic DTW algorithm is based on dynamic programming technique, which stores the previous computed results in a matrix so that the later calculation can use these results directly without re-computing them. Each cell in the matrix corresponds to an alignment of a point in both time series in each dimension - e.g cell (i, j) corresponds to the alignment of *i*th and *j*th time stamp in each time series. The value represents the DTW distance between the time series starting from the first data point up to the corresponding alignment point.

Therefore, given two time series, $X = \{x_1, x_2, \dots, x_m\}$ of length m and $Y = \{y_1, y_2, \dots, y_n\}$ of length n. The DTW distance

Big Data Mining and Analytics, September 2018, 1(1): 000-000

is calculated by following equations:

$$d(0,0) = 0$$

$$d(i,0) = d(0,j) = \infty$$

$$d(i,j) = || x_i - y_j || + min \begin{cases} d(x_{i-1}, y_j) & (1) \\ d(x_i, y_{j-1}) \\ d(x_{i-1}, y_{j-1}) \end{cases}$$

$$D(X,Y) = d(m,n)$$

Where $i = 1, 2, \dots, m$ and $j = 1, 2, \dots, n$ and D(X, Y) is the DTW distance between X and Y.

 $|| x_i - y_j ||$ is the distance between x_i and y_j , which can be $(x_i - y_j)^2$ or simply $| x_i - y_j |$, or other possible measures.

3.3 Distance Measure

Given two time series X and Y, the distance between any two points x_i and y_j from these two time series can be computed by *cost* function. There are many choice of the *cost* function, the most intuitive way is to use Manhattan distance, which is the absolute difference between x_i and y_j :

$$Manhattan_{distance}(x_i, y_j) = abs(x_i - y_i) \tag{2}$$

Euclidean distance is a well known method in measuring distance, which is the root of square difference:

$$Euclidean_distance(x_i, y_j) = \sqrt{(x_i - y_j)^2}$$
(3)

Another widely used measurement is the normalised distance, which measures the difference between the two points in the sequence with regard to the range of the sequences. That is, the difference between the maximum and minimum value in the sequences: $R_x = |X|, R_y = |Y|$. The normalised Manhattan distance can be computed with the following formula:

$$Manhattan_distance_norm(x_i, y_j) = \frac{abs(x_i - y_i) \times 2}{R_x + R_y}$$
(4)

3.4 GPGPU and CUDA

Using GPU as a general computing unit has attracted considerable attention. GPU provides massive parallel processing power. As the host for the GPU device, CPU organizes and invokes the kernel functions that execute on GPU. As shown in Figure 1(a), a GPU device consists of a number of streaming multiprocessors (SM), each comprising simple processing engines, called CUDA cores in the NVIDIA terminology [29]. Each SM has its own shared memory, which is equally accessible by all CUDA cores in the SM. At any given cycle, the CUDA cores in a SM execute the same instruction on different data items. SMs communicate with each other through the global memory of GPU.

From the programmers' perspective, the CUDA model [29] is a collection of threads running in



(a) The simple Cuda Hardware (b) The simple Cuda programming architecture mode

Fig. 1 CUDA programming Model

parallel. A warp is a collection of threads that can run simultaneously on a streaming multiprocessor (SM). The warp size is fixed for a specific GPU. The programmer decides the number of blocks and threads to be executed. If the number of threads is more than the warp size, they are time-shared internally on the SM. A collection of threads (called a block) runs on a multiprocessor at a given time. Multiple blocks can be assigned to a single multiprocessor and their execution is time-shared. A single execution generates a number of blocks. A collection of all blocks in a single execution is called a grid (Figure 1(b)). All threads of a block executing on a single multiprocessor divide its resources equally amongst themselves. Each thread and block is given a unique ID that can be accessed within the thread during its execution. Each thread executes a single instruction set called the kernel.

4 Algorithm for finding all similar subsequences

In this section, we propose a novel subsequence pattern mining algorithm. The objective of our algorithm is to discover all similar subsequences between two time series. Our proposed algorithm is based on the classical dynamic time warping algorithm and is adapted so that it can capture all subsequence patterns. We first introduce an intuitive strategy to find all similar subsequences. Then we propose two novel pattern mining algorithms.

4.1 Naive Subsequence Matching

The naive approach is to break the time series X into all possible subsequences, and compare them with all the subsequences of Y using the DTW algorithm and select the best matches based on the defined matching condition. That is, for all $X_s = (x_i, x_{i+1}, \ldots, x_{i+k})$, where $1 \le i \le n - k$, compute the DTW distance against all $Y_s = (y_j, y_{j+1}, \ldots, y_{j+l})$, where $1 \le j \le$ m-l, n and m are the length of X and Y respectively.

This approach will compute m DTW matrix for every timestamp in X, this is because it needs to examine all possibilities. Therefore, the complexity of this method is in quadratic time: $O((1+2+\cdots+n)(1+2+\cdots+m)) = O(n^2m^2)$.

Another brute force method is to use the shotgun window and try to match every possible window between the two sequences. For each window in the first sequence, compute DTW for all windows in the second sequence, and select the pairs with the best match, that is, with the lowest DTW cost. Figure 2 illustrates this approach. The complexity of the shotgun window approach is $O(m^2n^2/w^2)$, which only reduces the complexity by w^2 times. These naive approaches are not suitable for large-scale datasets, as the time complexity increases quadratically.



Fig. 2 The brute force shotgun windows approach. In these figures, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point.

Both algorithms presented in this section need to examine all possible combinations between different time series. Hence, the time complexity of both algorithms is high, which motivates us to design more efficient algorithms to find all similar patterns between two time series.

4.2 Proposed algorithm

The intuitive algorithms work by applying the DTW algorithm on all subsequences in two time series. Recall that the DTW algorithm works by adding the distance between two data points from two time series together. Therefore, we argue that if the distance between two

Big Data Mining and Analytics, September 2018, 1(1): 000-000

data points is large, it is less likely that it will lead to a similar pattern.

To verify our hypothesis, we conducted the following experiment. For two time series X and Y, we use the cost function to compute the distance between every data point and store the results in a matrix. We call this matrix the cross-distance matrix between two sequences, the value stored in each cell represents the distance between data points x_i and y_j . We visualise this matrix in Figure 3. In this figure, the darker area represents small values, and they correspond to the data points that have a shorter distance. From Figure 4, we can see that the alignments lie in the low-cost areas in the matrix, where the colours are the darkest in the visualisation.



Fig. 3 Normalised Manhattan Cost Matrix. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the Distance/Cost matrix represent the distance between two corresponding data points.



Fig. 4 Annotated Cost Matrix. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the Distance/Cost matrix represent the distance between two corresponding data points.

Based on the finding above, we modify the classic DTW algorithm so that it will only search for similar subsequences from the low-cost values cells.

In this algorithm, we first determine the low-cost values of the input time series. To achieve this, we introduce the cost threshold t, which is defined as the minimum average cost for each matching data points. This threshold t is used to determine whether to start or continue a DTW calculation. The procedure should accept the alignment (i, j) $(1 \le i \le N \text{ and } 1 \le j \le M)$ if $\frac{dtwDistance}{dtwLength} < t$, where:

$$dtwDistance = \sum_{i,j \in warppingPath} cost(x_i, y_j)$$
(5)

and dtwLength is the length of the warpping path.

If the DTW distance is smaller than this threshold, the path is considered as a possible match and will be included in the further computation.

To avoid over-warping in the warping path during the searching, the window constraint is imposed on the computation. In this method, we define an offset threshold o to control the length of the warping window, as illustrated in Figure 5. In this figure, the warping window is defined by the red region, and the threshold o is the length of the warping window offset from the starting position. With the offset threshold, the computation of DTW will stop when the computation reaches the warping window boundary, and therefore, the warping path will not exceed the length of the warping window.



Fig. 5 Offset Threshold and Warping Window

In some cases, the matching subsequences founded by the DTW algorithm may be too short to be considered as a valid match. To tackle this problem, we define the window threshold w to filter out the subsequences that are too short. The window threshold w is defined as the minimum alignment length of a sequence. As shown in Figure 6, suppose the red line

is a warping path, the length of the warping path must be longer than w for the subsequence to be considered.



Fig. 6 Window Threshold: Warping Path Length

4.3 Formal Definition of the algorithm

Based on the discussion in the previous sections, we first summarise the problem as the following:

Given sequence X = (x1, x2, ..., xn) of length n and sequence Y = (y1, y2, ..., ym) of length m, the algorithm searches for all subsequence $S = s_i, s_{i+1}, ..., s_{i+k}$, where S is a set of all subsequence of both X and Y that satisfy the following conditions:

- (1) The cost of the alignment per data point is less than t, that is, $\frac{d}{l} \leq t$. Where : d is total cost of the alignment, i.e. the sum of the cost of every alignment point in the warping path of the matching subsequence; l is the length of the warping path, i.e. the total number of the alignment points in the warping path.
- (2) The offset of the warping path from the perfect diagonal line must not exceed *o* data points, as explained in subsection 4.2.
- (3) The length of the warping path must be longer than the window threshold w, as explained in subsection 4.2. If the found match is too short, i.e. smaller than w, this match will not be included in the final result.

The proposed algorithm is based on the dynamic programming, which updates a matrix of size $m \times n$, where m and n are the length of two time serials X and Y respectively. In our algorithm design, we first create the following matrix to store the intermediate and final results:

• D: This matrix is used to hold the cumulative time warping cost, which takes O(mn) space.

- C: This matrix is used to store the distance between two data points. It is worth noting that this matrix is optional, since the cost between two data points can be computed on the fly while updating D.
- L: This matrix stores the current length of the warping path of the corresponding alignment. We increment the length of warping path while calculating the cumulated cost and finding the matched subsequences. We use the value stored in this matrix to determine if the match can be included in the final results.
- *R*: This matrix is the region-marking matrix that stores the position of the start point and end point of the matched subsequences, which will be used in the lookup and the trace-back processes.
- *P*: This matrix is used to store the direction of previous alignment point of the warping path, which is used in trace back to find the warping path.

The proposed algorithm is performed in three steps: initialisation, update the matrix, and find the path(trace back), which are detailed as follows.

- (1) Initialisation: The algorithm initialises all the matrices to zero and initialises cumulative cost matrix D to infinity. Therefore, during the update, the infinity will be ignored in the minimum comparison. If required, the cost matrix C can be computed.
- (2) Subsequence DTW calculation : This step performs the modified DTW algorithm. The algorithm starts from the top left position of the matrices, and iterates these matrices in the diagonal order.

If the current cell in the matrix is not part of any warping path and the cost of the current alignment is smaller than the cost threshold t, we mark this position as the start point in the region marking matrix R. Then we set the length of the warping path from the current cell in length matrix L as 1, update the cumulative cost to the current cost, and update other matrices accordingly. Otherwise, if the cost of current alignment is not smaller than threshold t, then we move on to the next position. If the cell being examined belongs to a warping path, then we select the cell with the smallest cumulative cost from the warping path. Next, we compute the average cost by adding the cost of the current position to the cumulative cost and divide this value by the total length. If the sum of the average cost and the cost of the current cell are smaller than threshold t, we include this cell as part of the matching subsequence from the cell with the smallest cumulative cost in the warping path, and update L, D, R, P respectively.

If all the elements in the warping path are not part of any matching subsequences, we first determine if the cost of this cell is smaller than t. If it is, then we start a new subsequence from the current cell and update all the matrices accordingly.

(3) Finding Paths: in this step the algorithm simply uses matrix R and L to find all the matched subsequences. The algorithm also uses threshold w to check if the length of the warping path is valid, then use P to trace back the warping path and return the results.

We outlined the above procedure in algorithms 1 and 2.

4.3.1 Further improvement

The algorithm described in the last section uses the minimum cumulative cost in matrix D to select the best cell for the DTW path. This approach requires the computation of the cumulative cost and the extra memory to store the results. Therefore, in this section, we propose a greedy algorithm to overcome this problem and further improve the performance.

The design of the greedy algorithm is based on the Markov assumption. That is, instead of using the cumulative cost, we only select the cell with the minimum cost from the warping path, and assume that the path from this cell will lead to a path with the lowest cumulative cost. With this assumption in mind, we change line 10 and line 12 in Algorithm 1 to Equation 6 and 7, respectively, and remove lines 16 and 22.

$$minpre = min(C[i, j-1], C[i-1, j], C[i-1, j-1]);$$
(6)

$$mini, minj = argmin([C[i, j - 1)], C[i - 1, j], C[i - 1, j - 1]);$$
 (7)

However, compared to the original design, the greedy algorithm does not achieve significant improvement in terms of performance and space. This is due

Procedure 1: DTW-based Subsequence Pattern Mining					
Input: Time series X and Y, threshold t and o Output: Matrices with subsequence matching info: D, L, R, P					
1 Initialisation: D to inf;					
2 L, R and P to 0;					
3 calculate cost matrix C ;					
/* begin dynamic programming	*/				
4 for $i = 1$: length(X) do					
5 for <i>j</i> = 1 : <i>length</i> (<i>Y</i>) do					
6 if $i == 1$ and $j == 1$ then					
7 $minpre = 0;$					
$ s \qquad \qquad mini = i; minj = j; $					
9 else $\min(D[i 1, i 1], D[i 1, i], D[i 1, i])$	<u>, 11)</u> ,				
$\begin{array}{c} minpre = (minpre = - inf) 2 0 \cdot min \\ minpre = (minpre$	7-1]),				
// set back to 0 if all previous	cells are				
still in init state					
12 mini, minj = argmin($D[i-1, j-1], D[i-1, j], D$	[i, j-1]);				
13 end					
$14 \qquad dtwm = (minpre + C[i, j])/(L[mini, minimum content in the c$	nj] + 1);				
if $dtwm \leq t$ and $L[mini, minj] == 0$ then					
/* start a new region	*/				
16 $D[i, j] = C[i, j];$ // update dt	w distance				
$\begin{array}{c c} 17 \\ D \\ $	ath length				
$\begin{array}{c c} & & \\ & &$	rt and end				
19 else if $dtwm \leq t$ then					
/* continue from previous region	*/				
20 (si, sj, li, lj) = R[mini, minj]; // find si,	ji as the				
region start cell					
// if current not diverge too muc	ch from				
offset					
21 if $abs((i-si) - (j-sj)) < o$ then					
$\begin{array}{c c} D[i, j] = minpre + C[i, j]; // \\ distance \end{array}$	update dtw				
23 $L[i, j] = L[mini, mini] + 1$:	// update				
path length	.,				
24 $R[i, j] = (si, sj, 0, 0);$ // up	date start				
cell					
25 $P[i, j] = (mini, minj); // u$	pdate path				
26 if $i > li$ and $j > lj$ then	beet out				
P_{i}	inest away				
$\begin{array}{c c} 2i \\ 28 \end{array} \qquad $					
29 end					
30 end					
ai end					

Procedure 2: Find Paths

I	Input: Matrices with subsequence matching info: L, R, P, threshold o								
C	Output: Matrices Path marking : OP								
1 Initialisation: OP to 0;									
2 fo	2 for $i = 1$: length(X) do								
3	3 for $j = 1$: length(Y) do								
	//	' if this is the start							
4	if	L[i, j] == 1 then							
5		_,_, $li,lj=R[i,j];$ // find end position							
6		if $L[li, lj] > w$ then // if path length > w							
7	while $li > si$ and $lj > sj$ do								
8	mark $OP[li, lj]$ as matching sequence path;								
9		(li, lj) = P[li, lj]; // get previous							
		step							
10		end							
11		mark $OP[li, lj]$ as matching sequence path;							
12		end							
13	en	d							
14	end								
15 e	nd								

to the following reasons. First, since the original algorithm is based on dynamic programming, in every iteration the cumulative cost is computed based on the result produced in the last step. Therefore, computing the cumulative cost is not an expensive operation. Second, although the greedy algorithm does not require matrix D to store the cumulative cost, it still uses all the rest matrices in computation. Hence, the space complexity is not reduced significantly. In addition, the greedy algorithm cannot find the optimal path because it only uses the local optimal results in finding the subsequence. However, in the DTW algorithm, the local optimal result may not lead to the global optimal solution.

5 GPU Acceleration

In this section, we discuss how to use CUDA to parallelize the algorithm presented in last section. We first introduce the parallelisation strategy used in our design, and then present the CUDA implementation. Further, we present our method to optimise this GPU implementation for the GPU memory hierarchy.

5.1 Data Dependencies

Recall that our algorithm uses the DTW algorithm as core to search for similar patterns. In DTW algorithm, however, the computation of each iteration depends on the result from the last step, which will limit the parallelisation degree during the computation. Therefore, we first analyze the data dependencies in our algorithm and then design a strategy to improve the parallelisation degree.

We show the data dependencies of the DTW computation in 7(a). In the matrix, each cell represents the DTW distance between two data points. The computation of each cell depends on the results from left, top, and top-left cells. That is, for cell (i, j) in the matrix D, the value of D[i, j] depends on the values in D[i, j-1], D[i-1, j], and D[i-1, j-1].

Therefore, for any cell (i, j) in the matrix, all cells between (0, 0) and (i, j) need to be computed before calculating the value of position (i, j). In the classic DTW algorithm, this computation is carried out in the row or column order. The cells in the same row or column are computed in sequence. On the other hand, there is no data dependency between the diagonal cells, as illustrated in 7(b). The cells connected by a red stripe have no data dependency. Therefore, these cells can be computed in parallel. The data dependency exist



Fig. 7 Matrix Update Order

between two neighbouring red strips. Therefore, we can only compute the data within one red strip at time. We illustrate this in 7(b) with grey arrow. The arrow indicates the data dependency and compute direction.

Based on the above discussion, we can summarise the parallel strategy as follows. The computation starts with the first cell, then updates the cells in the closest diagonal lines (the red stripe in 7(b)). Since these cells only depend on the first cell, they can be computed in parallel. Then the computation moves to the next diagonal line. By updating the matrix in this way, it is guaranteed that the data dependency is not broken and the reasonable degree of parallelism can be achieved.

5.2 Implementation of Parallelisation strategy

The previous section outlines a strategy to update the matrix in diagonal order. In this section, we describe how to implement this strategy in CUDA.

To implement this strategy on GPU, we need to consider how to map the CUDA threads to the data that need to be processed. Recall that the parallel diagonal computation model solves the data dependency issue by computing the matrix cells in diagonal order. Therefore, for each diagonal line, we can assign one thread to a cell within this line, and compute these cells in parallel.

An intuitive method to implement this strategy is to launch one kernel with enough threads, the number of threads within this kernel is larger or equal to the longest diagonal. In this design, the computation forms a loop within the kernel and each iteration processes one diagonal line. By implementing the algorithm in this way, we only need to launch one kernel, and therefore, reduce the overhead of multiple kernel launch.

In this implementation, we assign each column of the matrix to a thread first. Due to the data dependency, these threads are not processing each column simultaneously. Instead, each thread except



Fig. 8 Subsequence DTW Matrix Update Threading

the first one starts the processing when the previous thread finishes processing one cell in the matrix. By arranging the computation in this way, we can ensure that the data dependency is preserved and the data in the same diagonal line can be processed in parallel. This is because when the first thread processes the first cell in the matrix, the rest threads are waiting for this thread to finish. Once this computation is finished, the thread moves to the next item in the column, and its consecutive thread can start processing the first cell in the column assigned to it. These two threads can run in parallel because the cells involved in the computation can form a diagonal line in the matrix, and therefore, there are no dependency between them. However, as we mentioned in the last section, the results of the cells within a diagonal line depend on the cells in the previous diagonal line. Therefore, to enforce this dependency, we need to synchronise all threads involved in the computation along the diagonal line. We illustrate this design in Figure 8. In this Figure, the orange area represents the matrix that needs to be processed, the threads are represented by the red arrow, and the synchronisation point is labelled by the red diagonal line.

From this Figure, we can see that the computation

Big Data Mining and Analytics, September 2018, 1(1): 000-000

forms a parallelogram. The number of threads needed in the computation equals to the height of this parallelogram, which is ny. And the number of iterations required equals to the base of the parallelogram, which is nx + ny - 1 number of rows. Therefore, the time complexity of this implementation is: O(nx + ny - 1), which is a significant reduction from $O(nx \times ny)$.

5.3 Memory Access Optimisation

Coalesced access data in the global memory is critical to the performance of a GPU application. Non-coalesced memory access will lead to poor performance. In this section, we present how to enforce the coalesced memory access by enabling the consecutive threads to access consecutive memory address in our implementation.

Recall that our initial implementation processes the data in the diagonal order. However, the data is stored as a matrix in the memory, and the elements on one diagonal line are not stored consecutively in the memory. Therefore, to achieve the coalesced access to the global memory, the data stored in the memory need to be rearranged so that the elements on the same diagonal line are stored consecutively in the GPU memory.

Intuitively, rearranging the matrix can be achieved by reallocating another 2D array. In this new array, the length of x-dimension is the length of the longest diagonal line in the original matrix, while the ydimension is the number of the diagonal lines in the original matrix. In the new array, the diagonal lines from the original matrix are stored consecutively in the memory. We illustrate the array in Figure 9.

								0,0						
								1,1	0,1					
								2,1	1,2	0,2				
								3,1	2,2	1,3	0,3			
								4,1	3,2	2,3	1,4	0,4		
								5,1	4,2	3,3	2,4	1,5	0,5	
								6,1	5,2	4,3	3,4	2,5	1,6	0,6
							,	7,1	6,2	5,3	4,4	3,5	2,6	1,7
0,0	0,1	0,2	0,3	0,4	0,5	0,6		8,1	7,2	6,3	5,4	4,5	3,6	2,7
1,1	1,2	1,3	1,4	1,5	1,6	1,7			7.0					
2,1	2,2	2,3	2,4	2,5	2,6	2,7		8,2	7,3	6,4	5,5	4,6	3,7	
3,1	3,2	3,3	3,4	3,5	3,6	3,7		8,3	7,4	6,5	5,6	4,7		
4,1	4,2	4,3	4,4	4,5	4,6	4,7		8,4	7,5	6,6	5,7			
5,1	5,2	5,3	5,4	5,5	5,6	5,7		0 5	76	67				<u> </u>
6,1	6,2	6,3	6,4	6,5	6,6	6,7	1	8,5	7,0	0,7				
7,1	7,2	7,3	7,4	7,5	7,6	7,7	1	8,6	7,7					
8,1	8,2	8,3	8,4	8,5	8,6	8,7	1	8,7						
(a) Original Matrix								(b)	Dia	gona	l Mat	rix		

Fig. 9 Rearranging the matrix

However, as shown in the Figure, storing the diagonal line into the new matrix leads to the waste of memory space. To address this problem, instead of storing the diagonal line into a 2D array, we use a flat array to store the data from the diagonal line consecutively.

In addition, when calculating the warping boundary for the offset threshold, it is necessary to know the row and column position of each element. Therefore, we create another matrix, namely index matrix I, to store the index of the flat array in the corresponding row and column position.

The array indices can be determined through mathematical derivation, and be calculated in the initialisation step in parallel. In the next section, we present how to derive these index numbers in the matrix.

5.3.1 Calculating diagonal index

We draw some small matrix and fill in the index number to study the relationship between the array index and the row and column (i, j) in the matrix. The indices in the upper left corner and bottom-right corner in the matrix should be calculated differently. This is due to the following reason. If the index is computed from left to right, then the upper left part starts from the left most column, while the bottom right part starts from the bottom most row. As shown in the Figure 10 and Figure 11, these figures outlined two situations of the index number. The matrix represents the alignment of sequence X and Y, where the size of X is n_x and size of Y is n_y . The matrix size is $n_x \times n_y$; the row index is represented by i while the column index is represented by j, as shown in the figure. The red arrows in the matrix represents the direction of the increment of the array index.

Top-left:

As illustrated in Figure 10, when $j < n_x - i$, the array index falls into the top left part in the matrix. We denote the index of an element in the left most column as S_i , where *i* is the row index of the element. As shown in Figure 10(b), the number increases along the red arrow from the left most column. E.g. $index(0, 1) = S_1 + 1$; $index(1, 3) = S_4 + 3$.

For an arbitrary element residing in the top left part of the matrix (i.e. $j < n_x - i$), if we denote the start position of the diagonal line that this element belongs to as $S_{i_{start}}$, then we can compute its index by the following formula:

$$index(i,j) = S_{i_{start}} + j \tag{8}$$



Fig. 10 Optimised memory access index - top left

Because for any arbitrary element it forms a right triangle with x and y-axes, the position of i_{start} can be found by counting the number of steps needed for this arbitrary element to reach the leftmost column:

$$i_{start} = i + j \tag{9}$$

Since the number of elements in each diagonal line increases in each line, we can use the arithmetic sequence to calculate the $S_{i_{start}}$. However, as illustrated in Figure 10(c) and (d), when $n_x > n_y$, the computation has to be considered carefully. Since when the length of *x*-axis is greater than that of the y-axis, the arithmetic sequence stops at n_y and increases by n_y in each row, which can be formally defined as:

$$i_{start} > n_y \Rightarrow i+j > n_y$$
(10)

Hence, we separate the computation in following two cases:

• Case: $i + j < n_y$

Recall that the sum of an arithmetic sequence can be calculated by:

$$S_n = \frac{n(a_1 + a_n)}{2}$$

where :
 S_n : sum of *n* terms;

 a_1 : first term;

 a_n : n-th term

As illustrated in Figure 10(a) and (b) we get:

 S_i

$$_{start} = 0 + 1 + 2 + \dots + i_{start}$$
$$= \frac{i_{start} \times (i_{start} + 1)}{2}$$
(11)

Substituting Equation 11 with Equation 8 and Equation 9:

$$\begin{split} index(i,j) &= S_{i_{start}} + j \\ &= \frac{(i+j) \times (i+j+1)}{2} + j \end{split}$$

• Case: $i + j \ge n_y$

$$S_{i_{start}} = 0 + 1 + 2 + \dots + (n_y - 1) + (i_{start} - n_y + 1) \times n_y$$
$$= \frac{(n_y - 1) \times n_y}{2} + (i_{start} - n_y + 1) \times n_y$$
(12)

Substituting Equation 12 with Equation 8 and Equation 9:

$$\begin{split} index(i,j) &= S_{i_{start}} + j \\ &= \frac{(ny-1) \times ny}{2} + \\ &\quad (i+j-ny+1) \times ny + j \end{split}$$

Bottom-right:



Fig. 11 Optimised memory access index - bottom right

We now considering the case when $j \ge n_x - i$, i.e when the position falls into the bottom-right part of the matrix as shown in Figure 11. Similar to the topleft case, two cases need to be considered, as shown in Figure 11(a),(b) and (c),(d).

We denote the cells (i, j) in the bottommost row (i.e, where the red arrows start from) as $T_{j_{start}}$. The index(i, j) equals to $T_{j_{start}}$ adding the row index counting from the bottom (called the reverse row index). We denote the reverse row index by i', which can be computed by the formula:

$$i' = n_x - 1 - i \tag{13}$$

Hence, we can calculate the index(i, j) by:

$$\begin{aligned} udex(i,j) &= T_{j_{start}} + i' \\ &= T_{j_{start}} + n_x - 1 - i \end{aligned} \tag{14}$$

Similar to the Top-left case, the j_{start} is the number of steps required to reach the left most column from the current row:

$$j_{start} = j - i'$$

= $j - n_x + 1 + i$ (15)

The arithmetic sequence is used to calculate the $T_{j_{start}}$. In this case, the sequence is in the decreasing order from n_y to $n_y - 1 - j_{start}$. Hence, the arithmetic sequence starts at the point when:

$$j_{start} > n_y - n_x$$

$$\Rightarrow j - n_x + 1 + i > n_y - n_x$$
(16)

When $n_y > n_x$, the computation needs to be addressed differently. Hence, we handle the calculation in following 3 cases:

• Case: $n_y \leq n_x$

i

$$T_{j_{start}} = index(n_x - 1, 0) + (n_y + (n_y - 1) + \dots + (n_y - 1 - j_{start}))$$

= $index(n_x - 1, 0) + (17)$
 $\frac{(n_y + n_y - 1 - j_{start}) \times (j_{start})}{2}$

Substituting Equation 14 and Equation 15:

$$ndex(i,j) = index(n_x - 1, 0) + \frac{(2n_y - j + n_x - i) \times (j - n_x + 1 + i)}{2} + n_x - 1 - i$$
(18)

• Case: $n_y > n_x$ and $j_{start} \le n_y - n_x$

In this case, the $T_{j_{start}}$ is defined as:

$$T_{j_{start}} = index(n_x - 1, 0) + n_x \times j_{start}$$
(19)

By substituting Equation 14 and Equation 15, we can compute the index(i, j) as:

$$index(i, j) = T_{j_{start}} + n_x - 1 - i$$

= $index(n_x - 1, 0) +$
 $nx \times (j - n_x + 1 + i) + n_x - 1 - i$ (20)

• Case: $n_y > n_x$ and $j_{start} > n_y - n_x$ In this case, we compute the $T_{j_{start}}$ with the following formula:

$$T_{j_{start}} = index(n_x - 1, 0) + nx * (n_y - n_x) + (n_x + (nx - 1) + \dots + (n_y - 1 - j_{start})) = index(n_x - 1, 0) + n_x * (n_y - n_x) + (n_x + n_y - 1 - j_{start}) \times (j_{start} - (n_y - n_x))) 2 (21)$$

Again, substituting $T_{j_{start}}$ into Equation 14 and Equation 15 we get:

$$index(i, j) = T_{j_{start}} + n_x - 1 - i$$

= $index(n_x - 1, 0) + n_x * (n_y - n_x)$
+ $\frac{(n_y - j + i) \times (j + 1 + i - n_y)}{2}$ (22)
+ $n_x - 1 - i$

6 Evaluation

To evaluate the effectiveness and efficiency of the proposed pattern mining algorithm, we have conducted extensive experiments with different time series dataset drawn from real life. We first apply the proposed algorithm on different datasets to evaluate the ability of the proposed pattern mining method in capturing the repeated patterns between two time series. Then we evaluate the performance of the serial implementation and the parallel implementation. Finally, we evaluate our proposed technique for memory optimisation.

6.1 Preprocessing and parameters setting

We first evaluate the effectiveness of our proposed algorithm. In our experiments, the data preprocessing is needed before the pattern mining process. The preprocessing step includes removing trend, using Haar-wavelet-transfrom to smooth and reduce the complexity of time series with much noise and fluctuation. Preprocessing is beneficial to pattern mining, for example, the preprocessing steps such as Haar-wavelet-transform can reduce the size of the data, and hence will reduce the computation time.

In our experiments, we use the normalised cost threshold, offset threshold and window threshold. Hence, the same parameters can be applied on different datasets. In normalised cost threshold, we use the normalised cost function described in subsection 3.3. Therefore, the cost are in the range between 0 to 1. The time-series data should also be normalised in the preprocessing step if they are in similar ranges. In normalised offset and window threshold, we represent the offset/window as a factor of the average length of the time series instead of actual size. Therefore, the threshold parameters are independent on the dataset.

6.2 Subsequence Pattern Mining

A number of experiments have been conducted to evaluate the quality of proposed DTW-based pattern mining algorithm. In this section, we evaluate the algorithm with different datasets. The dataset used in these experiments are from the real world, synthetic data and random data.

6.2.1 Experiments with internet traffic data

Our first experiment is conducted by applying our algorithm on the datasets containing the same subsequence with different order. In this experiment, we use a subset of internet traffic data from The Time Series Data Library (TSDL) [30]. We crop a small piece of sequence, around 100 data points, as test dataset X. We also select another set of data with the similar length from the original dataset as dataset Y, and certain data in datasets X and Y are overlapped. In this experiment, we set the thresholds to the following values: t =0.13, o = 0.15, w = 0.2. The DTW cumulative cost matrix and paths for discovered subsequences are visualised in Figure 12 and Figure 13, respectively.



Fig. 12 Modified DTW method cumulative cost matrix. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the Distance/Cost matrix represent the distance between two corresponding data points.

The DTW cumulative cost matrix in Figure 12 shows that the DTW calculation starts from the start of potential matching subsequence, which is the low cost regions in the figure, the computation continues as long as the subsequence is still a eligible candidate. Figure 13 shows the results of trace back from the DTW cumulative cost matrix to report matching subsequence with the length more than the window threshold. These results demonstrate that the proposed algorithm is able



Fig. 13 Modified DTW method matched subsequence paths. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.

to capture the subsequence pattens between the time series. The exact subsequence is discovered, along with other similar subsequences.

This experiment shows that the modified DTW method picks up the local alignment regions in the matrix first, then calculates the possible subsequence alignments path. However, executing the algorithm with different distance function and thresholds will generate slightly different alignment results. Although the algorithm will find different alignments, these alignments will still correspond to the similar subsequence between two sequences.

6.2.2 Effect of different thresholds

The thresholds designed in the proposed algorithm can help the algorithm to discover the patterns and have impact on subsequence found by the algorithm. In this section, we study how the threshold values will affect the performance of proposed algorithm.

In these experiments, we use the same dataset from the previous section. This way, we can see the effect on the results with different thresholds. We first study the cost threshold t, which controls the toleration of the matching subsequence. In this experiment, we reduce the threshold t from 0.13 to 0.05. The results are shown in Figure 14.

The results shows that by reducing the cost threshold, the subsequences found by the algorithm lie in the lower cumulative cost in the path. Therefore, these subsequences are more similar. On the other hand,



Fig. 14 Result with lower cost threshold. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.

with larger value t, the algorithm will produce more matching sequences as it will consider more candidates.

We then study the effect of threshold o. This threshold controls how much the path can be diverged from the perfect matching sequence path (i.e a one to one correspondence in time warping path). If the offset threshold o is too big, the discovered path will include long tails as demonstrated in Figure 15, which uses a larger offset threshold of 1. On the other hand, if the offset is too small, the result may only include the straight path and they may be clustered together. Unlike other thresholds, the offset threshold o has a bigger influence on the quality of the result.

The window threshold is used to filter out the shorter sequences and only report the matching sequences that have the warping path longer than the threshold. In this experiment, we increase the window threshold from 0.2 to 0.3 and plot the results in Figure 16. From this figure, we can see that by increasing the window length, the results produced by the algorithm is less than the small threshold.

6.2.3 Experiments with the data containing spikes - WordsSynonyms

We now evaluate how the proposed algorithm performs when the datasets are not perfect. In real-world time series data, peaks or spikes can be observed frequently. Thus it is essential to know how our algorithm can handle this problem. In these experiments, we use the WordSynonyms time series



Fig. 15 Result with large offset threshold. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.



Fig. 16 Result with larger window threshold. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.

data from UCR Time Series Classification Archive [31], which contains the possible patterns but also contains the spikes in one of the sequences. In this dataset, the training and testing datasets are different but are from the same source.

We compare the training and testing data from this dataset to evaluate our pattern mining algorithm. Unlike the training data, the testing dataset contains spikes in it. In order to see how well our algorithm performs, we compare the results produced with the original dataset against the dataset with the spikes being removed. There are many different techniques to remove the spikes from the dataset. In our experiment, we choose Median filter to eliminate the spikes in time series dataset due to the popularity of this methods. The experiment results are shown in Figure 17 and Figure 18.



Fig. 17 Result of WordsSynonyms data. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.



Fig. 18 Result of WordsSynonyms data filtered. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.

From these figures, we can see that the matching subsequence results found in the original data are similar to the result obtained with the data in which the spikes are removed. These results prove that our method is resistant to spikes in the data and it will treat them as part of the pattern.

6.2.4 Experiments with the circular trend data

In this section we evaluate the performance of proposed algorithm with the data that have the circular patterns. In this experiment, we use the data with the repeating patterns and the increasing trend (the absolute value of the data increases over time) from TSDL [30]. The time series is tested against itself in this experiment to see how our method handles the type of time series data. We create another dataset by removing the trend in the data, and compare the results with the original dataset. As we can see from Figure 19 and Figure 20, our algorithm is able to is able to pick up all the circular patterns after removing the trend.



Fig. 19 Data with circular trend. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.

6.2.5 Experiments with noisy data - cluster load

We then evaluate our algorithm using the data with the noisy signals. In this experiment, we use two sets of trace data in clusters from [32]. The result alignments look very scattered without any preprocessing. To improve the result, we applied the Haar Wavelet transform [33] to the data sequence. Figure 21 and Figure 23 show the result before and after applying the Haar-Wavelet-transform to the two sets of data. Figure 22 and Figure 24 demonstrate that the paths of the patterns in the cost matrix. The results show that the patterns are discovered correctly.

6.3 Performance

In this section, we conduct the performance analysis on both serial and parallel implementation. In these



Fig. 20 Data with circular trend - removed trend. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.



Fig. 21 Haar wavelet transform on complicated time series. In these figures, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point.

experiments, we use a very large dataset collected by [34]. We test the performance by using different implementations to analyze a subset of the data with different lengths. The experiments are conducted on a machine with NVIDIA Tesla K40 GPU with 12GB memory. We use a subset from the dataset that can fit into the GPU memory.

6.3.1 Serial vs CUDA Implementation

We profile the implementation using the same data on the vertical and horizontal axis. The execution time is recorded with different lengths of the data. Figure 25 shows the execution time of the



Fig. 22 Modified DTW on transformed sequence. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.



Fig. 23 Haar wavelet transform on complicated time series. In these figures, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point.

serial implementation (Host_Matrix) and the CUDA implementation (CUDA_Matrix).

From the figure, we can see that the CUDA implementation achieves the faster execution time than the serial implementation.

We then compare the execution time between the CUDA implementation and the CUDA implementation with the optimised memory access.

It can be seen from Figure 26 that by optimising the memory access the optimised CUDA implementation has the shorter execution time. This result demonstrates the effectiveness of our technique in optimising the memory access.



Fig. 24 Modified DTW on transformed sequence. In two input data frames, the x-axis represents the length of input time series data (number of data points), the y-axis represents the value of each data point. The x and y-axes in the matrix represent the distance between two corresponding data points. The matching subsequences between two data frames are shown with the yellow line.



Fig. 25 Execution time of the serial and CUDA implementations. The x-axis represents the length of input time series data (number of data points), the y-axis represents the execution time in millisecond.



Fig. 26 Comparison of dtwm calculation time. The x-axis represents the length of input time series data (number of data points), the y-axis represents the execution time in millisecond.

7 Conclusion

In this paper, we proposed a novel algorithm to find all similar subsequences in the time series data. Our algorithm first identifies the possible start positions of similar alignments, and then uses the dynamic programming technique to expand the subsequences. Further, we parallelize the proposed algorithm using GPGPU and CUDA, and develop a optimization technique to improve the memory access in the GPU implementation. Extensive evaluation experiments have been conducted. The results show that our algorithm can discover similar subsequences between two time series and achieve much better performance with the GPU parallelisation.

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Big Data Mining and Analytics, September 2018, 1(1): 000-000

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Big Data Mining and Analytics, September 2018, 1(1): 000-000

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