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# An improved version of the runs algorithm based on Crochemore's partitioning algorithm * 

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#### Abstract

Though there are in theory linear-time algorithms for computing runs in strings, recently two of the authors implemented an $O(n \log n)$ algorithm to compute runs that was based on the Crochemore's partitioning repetitions algorithm. The algorithm preserved the running complexity of the underlying Crochemore's algorithm; however, the static memory requirement - already large at $14 n$ integers for a string of length $n$ - was increased significantly to $O(n \log n)$ integers. The purpose and advantage of this algorithm was its speed. In this paper we present a more advanced version of the extension of the Crochemore's algorithm for computing runs. This version in addition to maximal repetitions, computes runs and primitively rooted distinct squares. Its implementation completely does away with the extra memory required for the previous version and through some additional memory saving techniques, the overall memory need was reduced to $13 n$ integers.


Keywords: repetition, run, distinct squares, string, periodicity, suffix array, LCP array, Lempel-Ziv factorization

## 1 Introduction

Crochemore's repetitions algorithm, often also referred to as Crochemore's partitioning algorithm, was introduced in 1981 [4] and was the first $O(n \log n)$ - and hence optimal - algorithm to compute maximal repetitions in a string of length $n$. The big advantage of the algorithm was its independence on the size of the alphabet of the string. Its disadvantage was in the implementation, as the data structures required for keeping track of the refinement process and the gaps require a substantial storage - originally estimated in about $20 n$ of integers and a complex machinery to update and maintain them. In 2003, Franek, Smyth, and Xiao [9] implemented the algorithm using several memory saving techniques lowering the requirement to $14 n$ integers. An additional advantage of their implementation was that the memory was static, or to be more precise, allocated all at once at the outset of the algorithm as the working of the algorithm did not require any dynamic allocation or deallocation of memory. This approach led to an implementation with quite fast running times.

Since the advent of linear-time algorithms to compute suffix arrays [10, 12, 19], an avenue opened for true linear-time algorithms to compute runs. Such algorithms follow the general strategy of

[^0](a) compute suffix array using any of the linear-time algorithms, for instance [10, 12, 18-20],
(b) compute LCP (longest common prefix) array using any of the linear-time algorithms, for instance $[11,17]$
(c) compute Lempel-Ziv factorization using any of the linear-time algorithms, for instance $[3,5]$
(d) compute some runs that include all leftmost runs from the Lempel-Ziv factorization using Main's algorithm [15, 16]
(e) from the runs computed in (d), compute all runs using Kolpakov-Kucherov's approach $[13,14]$

The laborious and circuitous strategy for linear-time algorithms suggests that performance of such algorithms may not be satisfactory. Franek and Jiang [6, 7] extended the original Crochemore's repetitions algorithm to compute runs with a plan to benchmark the algorithm and compare it with any implementation of the linear-time algorithm for computing runs. Their implementation was based on Franek, Smyth, Xiao's implementation [9] for its optimized memory handling. The approach was quite straightforward: the maximal repetitions as reported were collected and consolidated into runs. This necessitated additional data structures of size $O(n \log n)$ integers. The program still exhibited fast running times, but the memory requirement was too substantial and required dynamic handling of memory during processing, which is quite a detriment to fast performance.

The reason to revisit the algorithm and modify it was to lower the memory requirement, eliminate the need of dynamic memory allocation and deallocation during processing, and prepare the stage for the parallelization. This report describes the new implementation that requires only a single allocation of $13 n$ integers at the outset of the algorithm, preserves all the advantages of the previous implementations, computes not only the maximal repetitions - as the original Crochemore's algorithm does, but also the runs - as the Franek and Jiang's implementation does, and in addition it computes the number of primitively rooted distinct squares. Moreover, the algorithm in this form is well-posed for parallelization in the sharedmemory model. We refer to this algorithm as FJW.

## 2 Preliminaries

For $e \geq 2$ and a non-empty string $w,(w w)^{e}$ is a repetition of power $e$ in a string $x$ if there are strings $u$ and $v$, possibly empty, so that $x=u(w w)^{e} v$. $w$ is referred to as the generator of the repetition, while the size of the generator is referred to as the period of the repetition. If $e=2$, we talk of a square. A string is primitive if it is not a repetition. A repetition is primitively rooted if its generator is primitive. A repetition $(w w)^{e}$ in $x=u(w w)^{e} v$ is maximal if $w$ is neither a suffix of $u$ nor a prefix of $v$. For a string $x=x[0 . . n-1]$, a repetition can be encoded as a triple $(s, p, e)$, where $s$ is the starting position of the repetition, $p$ is the period, and $e$ is the power.

A more succinct notion is that of a run. In a string $x=x[0 . . n-1]$ a quadruple $(s, p, e, t)$ encodes a run if
(a) for any $0 \leq i \leq t,(s+i, p, e)$ is a maximal repetition,
(b) either $s=0$ or $(s-1, p, 2)$ is not a square, i.e. the run cannot be extended to the left,

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(c) either $s+t=n-1$ or $(s+t+1, p, 2)$ is not a square, i.e. the run cannot be extended to the right,
(d) the generator $x[s . . s+p-1]$ is primitive.

The maximum number of maximal repetitions in a string of length $n$ is $O(n \log n)$, see [4]. On the other hand, the maximum number of runs is $\leq 1.029 n$, see [1]. In $[6,7]$, Franek and Jiang used Crochemore's repetitions algorithm to generate all maximal primitively rooted repetitions, collect them in a data structure of size $O(n \log n)$ and then in $O(n \log n)$ time process the collected repetitions and consolidate them into runs. Though the repetitions computed by Crochemore's algorithm are not in any particular order - except the fact that repetitions of the same period are computed at the same stage, a detailed examination of the gap function revealed that there is no need to collect the repetitions, that the runs can be directly inferred from the information provided by the gap function.

To be able to discuss the gap function and show how the runs can be directly inferred, we need to briefly discuss the mechanism of the Crochemore's repetitions algorithm.

## 3 Brief description of Crochemore's repetitions algorithm

In mathematical terms, the algorithm is simple and elegant and relies on the refinements of classes of equivalence of the positions of the input string $x=x[0 . . n-1]$. An equivalence $\approx_{k}$ is defined on the set of indices $\{0, . . n-1\}$ by $i_{1} \approx_{k} i_{2}$ if and only if $x\left[i_{1} . . i_{1}+k-1\right]=x\left[i_{2} . . i_{2}+k-1\right]$. In simple terms, two positions are $\approx_{k}$ equivalent, if the substrings of length $k$ starting at those two positions are the same. In all times, the algorithm maintains an ascending order of the indices in each class, though no particular order of the classes themselves.

At the first level, the algorithm computes by brute force the classes of equivalence $\approx_{1}$. These classes in fact represent all the positions with the same alphabet symbol. On each following level $k$, all classes of equivalence $\approx_{k}$ are computed. Note that each class from level $k-1$ is either preserved as a class on level $k$, or is partitioned into several disjoint classes which we will refer to as family. That is why the Crochemore's algorithm is also referred to as the partitioning algorithm. It is clear that once a class has size 1 , it cannot be partitioned any further. The processing ends when all classes are of size 1 .

The classes, indeed, contain all information of all possible repeats of substrings of $x$. It is straightforward to see that a primitively rooted square of period $p$ must be represented by two consecutive indices $i_{1}$ and $i_{2}$ in the same class of $\approx_{p}$ so that $\left|i_{1}-i_{2}\right|=p$.

The main complication of the algorithm lies in the process of refinements. If the refinements were carried out directly through references to the input string, the running complexity would be unacceptable $O\left(n^{2}\right)$. However, the refinement of the class on level $k$ can be carried out by using other classes on level $k$ which allows to discard the original string once the classes on the first level had been computed. This approach, though much better than the refinement through direct reference to the input string, would still lead to the running complexity of $O\left(n^{2}\right)$. If in each family we take a largest class by size and designate it big and all other as small, we can carry the full refinement of all the classes using just the small classes. Since any position can occur in at most $O(\log n)$ small classes, this approach gives the running complexity of $O(n \log n)$.

Not to destroy the $O(n \log n)$ complexity, we cannot afford to scan the classes when looking for squares and ultimately for maximal repetitions. Throughout the whole process of refinement, a function $\operatorname{Gap}(p)$ is maintained that gives a list of all indices that are exactly $p$ distance from its predecessor in the class, more precisely: when processing level $k$, if $i_{2} \in$ $\operatorname{Gap}(p)$, then $i_{1}=i_{2}-p$ is in the same class of equivalence $\approx_{k}$ as $i_{2}$ and these two indices are consecutive in the class. We will describe the Gap () function in more detail in the next section dealing with the implementation of the FJW algorithm.

## 4 Implementation of the FJW

We first describe the implementation and its data structures without any regard for the size of required memory. This leads to an implementation requiring $19 n$ of integers. Then we use several techniques to reduce the required memory to $13 n$ integers. We will present the data structures as static, but for practical reasons - we do not want to recompile the program each time a different string is to be processed, all the structures are allocated once at the outset of the program's processing. The structures are essentially arrays used to emulate doubly-linked lists, stacks, and queues.
The first seven arrays deal with classes:

1. An integer array $C \operatorname{Start}[0 . . n-1]$ stores the very first element of a class, i.e. $C S t a r t[i]=j$ means that the first element of class $i$ is $j$. This emulates a pointer to the beginning of a class.
2. An integer array $C E n d[0 . . n-1]$ stores the very last element of a class, i.e. $C E n d[i]=j$ means that the last element of class $i$ is $j$. This emulates the pointer to the end of a class.
3. An integer array $C N e x t[0 . . n-1]$ stores the next element in the class or null. Thus $C N e x t[i]=j$ indicates that $i$ and $j$ are in the same class and that $j$ is the next element after $i$, while $C N e x t[i]=$ null indicates the $i$ is the last element in the class. This emulates the forward links.
4. An integer array CPrev[0..n-1] stores the previous element in the class or null. Thus $C \operatorname{Prev}[i]=j$ indicates that $i$ and $j$ are in the same class and that $j$ is the element just before $i$, while $C \operatorname{Prev}[i]=$ null indicates the $i$ is the first element in the class. This emulates the backward links.
5. An integer array $C M e m b e r[0 . . n-1]$ stores the membership of each element, i.e. $C M e m b e r[i]=$ $j$ means that $i$ belongs to the class $j$.
6. An integer array CSize $[0 . . n-1]$ stores the sizes of classes, i.e. CSize $[i]=j$ means that class $i$ has size $j$.
7. An integer array CEmpty $[0 . . n-1]$ is used as a stack of empty classes to be used.

The following four arrays deal with families:

1. An integer array $F \operatorname{Start}[0 . . n-1]$ is used as a stack. $F \operatorname{Start}[i]=j$ thus means that class $j$ is the first class in the family $i$.
2. An integer array $F N e x t[0 . . n-1]$ emulates the forward links in a list of classes in a family.
3. An integer array $F \operatorname{Prev}[0 . . n-1]$ emulates the backward links.

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4. An integer array $F M e m b e r[0 . . n-1]$ stores the family membership, i.e. $F M e m b e r[i]=j$ means that class $i$ belongs to family $j$.

The following four arrays deal with the refinement process:

1. An integer array Refine $[0 . . n-1]$. Refine $[i]=j$ means that an element from class $i$ should be moved to class $j$.
2. An integer array $\operatorname{RStack}[0 . . n-1]$ is used as a stack. It is used to remember which items in Refine[] were occupied, so it can be cleared without any need to traverse the whole array Refine[] which would destroy the $O(n \log n)$ complexity.
3. An integer array $\operatorname{Sel}[]$ is used as a queue. It is the queue of all elements of all small classes.
4. An integer array $S c[]$ is used as a queue of small classes. $S c[i]=j$ indicates $j$ is the last element of a small class. Thus the information in $S e l[]$ and $S c[]$ implements a list of elements of small classes with indicators where one small class ends and the next small class starts.

The last four arrays implement the gap function:

1. An integer array $\operatorname{Gap}[0 . . n-1] . \operatorname{Gap}[i]=j$ indicates that the first element in the gap list for $i$ is $j$, i.e. $j$ 's predecessor in the class is $j-i$.
2. An integer array GMember $[0 . . n-1]$. GMember $[i]=j$ means that $i$ belongs to the gap list $j$.
3. An integer array $G N e x t[0 . . n-1]$ emulates the forward links in the gap lists.
4. An integer array GPrev[0..n-1] emulates the backward links in the gap lists.

The C++ code for this version is the file crochB.cpp and electronically available at [2].
In the next version, crochB1.cpp, also posted at [2], the array GMember[] is replaced by a function $G M e m b e r()$ and the memory requirement is reduced to $18 n$ integers. GMember () can be directly computed:
$\operatorname{GMember}(i)= \begin{cases}\text { null } & \text { if } i \text { is not member of any class, } \\ \text { null } & \text { if } i \text { is the first member of a class, } \\ i-C \operatorname{Prev}[i] & \text { otherwise. }\end{cases}$
Version crochB3.cpp reduces the memory requirement further to $17 n$ integers. Consider any family doubly-linked list, its beginning can be determined by two means: $F S t a r t[i]=j$ or FPrev $[j]=$ null. Thus, we can do away with $F M e m b e r[]$ array and replace it by a function that is utilizing the redundant space in FStart[] and FNext[]:
$F M \operatorname{mber}(i)= \begin{cases}F \operatorname{Start}[i] & \text { if the stack pointer is null, } \\ F N \operatorname{ext}[F \operatorname{Prev}[F \operatorname{Start}[i]]] & \text { if } i \leq \text { the stack pointer, } \\ F \operatorname{Start}[i] & \text { otherwise. }\end{cases}$
We also introduce a function $\operatorname{FEnd}()$ computed from FStart [] and FPrev[]: FEnd $(i)=$ FPrev[FStart $[i]]$.

In the next version, crochB4.cpp $C E m p t y[]$ and $S c[]$ are made to share the same memory segment, reducing the memory requirement to $16 n$ integers.

Version crochB5.cpp distributes CEnd[] and CSize[] over CStart, CNext, and CPrev, thus reducing the memory requirement further to $14 n$ integers. Therefore, $C E n d(i)=C \operatorname{Prev}[C S t a r t[i]]$ and $\operatorname{CSize}(i)=C N e x t[C \operatorname{Prev}[C \operatorname{Start}[i]]]$.

If we limit the maximal possible length of an input string from UNSIGNED LONG MAX to LONG MAX, which for a 32 -bit long it is $2,147,483,647$ and thus large enough, we can virtualize CMember[] over Gap[], GNext[], and GPrev[], reducing the memory requirement to $13 n$ integers. Thus, the function to set the value of $C M e m b e r(e)$ to $c$ :

```
if (Gap[e] == null || Gap[e] < 0)
    if (c == null)
            Gap[e] = null;
    else
            Gap[e] = 0-1-c;
else
    if (c == null)
        GNext[GPrev[Gap[e]]] = null;
    else
        GNext[GPrev[Gap[e]]] = 0-1-c;
```

and the function to get the value of $C$ Member $(e)$ :

```
if (Gap[e]==null)
    return null;
else
    if (Gap[e] < 0)
        return 0-1-Gap[e];
    else
        if (GNext[GPrev[Gap[e]]] == null)
            return null;
        else
            return 0-1-GNext[GPrev[Gap[e]]];
```

The version crochB7.cpp is just a polished version of crochB6.cpp with the additional features discussed in the next section.

## 5 The gap function and computations of distinct squares, maximal repetitions, and runs

Throughout the process of refinement, the gap function is maintained. In order to protect the running complexity of $O(n \log n)$, every time an element is removed from a class, the gap function is updated; and any time an element is added to a class, the gap function is updated again. When computing the next level from the current one, $\operatorname{Gap}[p]$ points to the first element whose immediate predecessor in its class is exactly at distance $p$, while $G N \operatorname{ext}[]$ and GPrev[] allow us to traverse the whole list in either direction and to update the list in

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constant time. Notice that if $G a p[p]=i$ and we are dealing with level $p$ of refinement, then there is a primitively rooted square starting at position $\operatorname{GPrev}[i]$ of period $p$.
Computing Distinct Squares - traceSquares() method
The gap function can be used to compute primitively rooted distinct squares. As we traverse the gap list, once we identify the first primitively rooted square in each class, we ignore the identification of the rest from the same class as they are all identical squares. We use Refine[] and RStack[] that are only needed during the refinement process as auxiliary data structures here to indicate the last class we already have a representative from in order not to get another representative from the same class. Note that the program can either output the number of distinct squares, the triples $(s, p, e)$ encoding the squares identified, or the squares as identified substrings of the input string - we refer to is as pretty print. To use pretty print, the string alphabet should be the lower case letters $a, b, \ldots$.
Computing Maximal Repetitions - traceMaxReps() method
For the maximal primitively rooted repetitions, again either their number can be output, the individual repetitions in their encoding into triples or pretty print can be used. The algorithm traverses the gap list, and for each entry it checks how far left and how far right it can extend the square. Thus, during the tracing at level $p$, all the individual squares identified are consolidated into maximal repetitions. A brief description on how the algorithm determines if the square can be extended to the left: the entry $i$ from the gap list Gap $[p]$ indicates that there is a primitively rooted square starting at position $i-p$. Then the algorithm checks if the square can be extended to the left - i.e. is there a square of period $p$ starting at position $i-2 p$ and determined by $i-p$. It is possible that the position $i-p$ is in the gap list further away. In order not to process the square starting at $i-2 p$ and determined by $i-p$, we again use Refine[] and Rstack[] to indicate that this entry has already been processed.
Computing Runs - traceRuns() method
The computation of runs is performed by TraceRuns(). The idea is very similar to that of tracing maximal repetitions: the identified primitively rooted squares are consolidated to runs. If you look at the leading square of a run $(s, p, e, t)$ that must be primitively rooted by definition, at every position $s+i, 0 \leq i \leq(e-2) * p+t$ there is a primitively rooted square. This fact is based on a simple observation that a rotation of a primitive string is also primitive:

Lemma 1. If a string $\boldsymbol{u}$ is primitive, then every rotation of $\boldsymbol{u}$ is also primitive.
Proof. It suffices to prove that if $\boldsymbol{u}[0 . . k-1]$ is primitive, so is $\boldsymbol{u}_{\boldsymbol{1}}=\boldsymbol{u}[k-1] \boldsymbol{u}[0 . . k-2]$.
Assume that $\boldsymbol{u}_{\mathbf{1}}$ is not primitive. Then $\boldsymbol{u}_{\mathbf{1}}=\left(v_{1} \ldots v_{r}\right)^{q}$ where $1 \leq r$ and $2 \leq q$. So, $\boldsymbol{u}_{\boldsymbol{1}}=v_{1} \ldots v_{r}\left(v_{1} \ldots v_{r}\right)^{q-2} v_{1} \ldots v_{r}=v_{1} \ldots v_{r-1}\left(v_{r} v_{1} \ldots v_{r-1}\right)^{q-2} v_{r} v_{1} \ldots v_{r}$. Thus, $\boldsymbol{u}=v_{r} v_{1} \ldots v_{r-1}\left(v_{r} v_{1} \ldots v_{r-1}\right)^{q-2} v_{r} v_{1} \ldots v_{r-1}=\left(v_{r} v_{1} \ldots v_{r-1}\right)^{q}$ implying that $\boldsymbol{u}$ is not primitive, a contradiction.

In the algorithm, we have to consolidate the run from all of the primitively rooted squares encoded in the gap function. Thus, having identified a square, not only we must check if it can be extended left or right as a repetition, we have to check if it can be shifted left or right. Again, we are using Refine[] and RStack[] as auxiliary data structures to indicate which of the elements of the gap list had been previously processed as the part of tracing, so we do not process them again.

## 6 Preparations for parallelization of the FJW

The process of refinement of a class by other classes is quite independent from any other refinements, and thus can be easily carried in parallel. In a shared memory model, such as a multi-core machine, all processors carrying on the individual refinements can access and use all the same data structures, provided a proper care for locking and unlocking is taken for the updates of these data structures. [8] discussed two different approaches to parallelize the refinement step of the Crochemore's algorithm. One of them is to assign each processor a single class to be refined by all the small classes, and all the classes are being refined in parallel. The advantage of this approach is that it no longer requires refinement structures of Refine[] and RStack[] as the serial version does because each processor only refines one class and thus only one destination class needs to be carried for each refinement by a particular small class at a time. Therefore, for this approach each processor just needs an extra integer to carry on the refinement. The disadvantage of this approach is that each processors may take almost as long as a single processor in the serial implementation, as it must traverse the entire small classes. Though there are some potential speed-ups based on the fact of knowing the smallest and the largest elements of the class to be refined, so that limits the elements of each small class that needs to be traversed. As the implementation of FJW where each element $e$ in the small classes is processed and element $e-1$ is refined from some classes, the second approach of the parallelization is to assign each processor a small class to be processed, and all the small classes being processed in parallel. This approach can dramatically improve the computation speed since there is no unnecessary work being done as the first approach does, and the elements in small classes only been traversed once in total. However, a separate copy of refinement structures of Refine[] and RStack[] is required for each processor since when two processors refining the same class by two different small classes, conflicts may occur and therefore would lead to incorrect results. [8] suggests some strategy to reduce this requirement though still fairly amount of extra memory is required and it involves the dynamic memory allocation which would slow down the computation. To overcome this issue, succinct data structure bit vectors can be used to avoid the extra memory allocation. Therefore, for both approaches discussed above, the data structures as used in the sequential algorithm can be be used as they are in the parallel version. The task of parallelization is thus reduced to the proper protection of the shared data structures during the updates.

## 7 Conclusion

We present a new implementation of an extension of the Crochemore's repetitions algorithm that computes primitively rooted distinct squares, primitively rooted maximal repetitions, or runs. The running complexity of the algorithm is preserved and thus is $O(n \log n)$ where $n$ is the length of the input string. In comparison to the previous implementation of the Crochemore's partitioning algorithm, the memory required is reduced to $13 n$ integers. In comparison to the previous implementation of an extension to compute runs, there is no additional memory requited and no dynamic allocation or deallocation of the memory during the processing as all the required memory is allocated once at the outset of the program.

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The resulting algorithm implemented in C++ is very fast and all the versions described in this paper can be downloaded from [2]. The future work will include the bench-marking of the algorithm and comparisons with other implementations of algorithms for computing of runs, especially the linear ones.

We discussed the issues of parallelization and concluded that this implementation can be parallelized in the shared memory model with very little effort and will be, thus, done in a short term future.

One of the author, C. Weng, is currently working on the implementation of a true lineartime algorithm for computing runs. Once the implementation is complete, the FJW algorithm will be bench marked and tested against the linear-time algorithm.

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