**IndexToolkit: an open source toolbox to index protein databases for high-throughput proteomics**

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**ABSTRACT**

Summary: A software package, IndexToolkit, aimed at overcoming the disadvantage of FASTA-format databases for frequent searching, is developed to utilize an indexing strategy to substantially accelerate sequence queries. IndexToolkit includes user-friendly tools and an Application Programming Interface (API) to facilitate indexing, storage and retrieval of protein sequence databases. As open source, it provides a sequence-retrieval developing framework, which is easily extensible for high-speed-request proteomic applications, such as database searching or modification discovering. We applied IndexToolkit to database searching engine pFind to demonstrate its effect. Experimental studies show that IndexToolkit is able to support significantly faster searches of protein database.

Availability: The IndexToolkit is free to use under the open source GNU GPL license. The source code and the compiled binary can be freely accessed through the website http://pfind.jdl.ac.cn/IndexToolkit. In this website, the more detailed information including screenshots and documentations for users and developers is also available.

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**INTRODUCTION**

Protein identification is a critical step in most high-throughput proteomics research. A huge number of tandem mass spectrometry (MS/MS) data needs to be searched in protein databases to identify the protein sequences ( Aebersold and Mann, 2003 ). Currently, most popular protein sequence databases, such as Swiss-Prot and IPI ( Kersey et al., 2004 ), are represented in the FASTA format, which is a flat text. However, it is inefficient to perform frequent searches directly on such text-based databases since in each search the whole database has to be scanned from the first entry to the last. In general, when a database is frequently searched but less often read into memory, acting with indexes. IndexToolkit also supports memory mapping to read the whole or a part of the index database into main memory.

Recently, a software package named DBToolkit ( Martens et al., 2005 ) is made available publicly to convert protein sequence database into peptide sequence database to enhance protein identification. However, the output of DB Toolkit is still a text file, and it does not support indexing. As far as we know, there is no open-source package to help experimenters index protein/peptide sequence database [note that the open source search engines X!Tandem ( Craig and Beavis, 2004 ) does not provide general-purpose indexing support. Although commercial search engines such as Mascot and SEQUEST create indexes for their own utilization, they do not provide indexing support for other researchers to develop their own search engine. Although various new peptide-scoring algorithms are being proposed, the lack in the support of efficient database indexes hampers them from practical application.

To solve this problem, we have developed an open source software called Index Toolkit to perform indexing for a protein sequence database. Index Toolkit can help researchers who develop search engines themselves to improve greatly the efficiency of high-throughput protein identification.

**2 DESCRIPTION OF INDEX TOOLKIT**

Here we briefly describe the implementation, usage and features of Index Toolkit. Index Toolkit consists of two parts: a set of API ( Application Programming Interface ) functions for application developers and user-friendly software tools for common users ( Fig. 1a ). API is the heart of Index Toolkit and is written in the C++ language with Standard Template Library ( STL ). The Index Toolkit API is open as a C++ class library, including a FASTA format parser, digestion and fragmentation simulator, index creator, index loader, index query and other sequence processing functions. Using the classes, developers can execute various operations interacting with indexes. Index Toolkit also supports memory mapping to read the whole or a part of the index database into main memory.

Tools in the Index Toolkit package are implemented to help users to manage the index more effectively, such as importing, setting, indexing and retrieving databases. All tools are developed based on the Index Toolkit API and their sources demonstrate how to use API. Users can realize their own powerful tools in two ways: directly use API or modify the source codes of Index Toolkit tools. As an open project, developers can easily add Index Toolkit into their projects and simultaneously provide a high-speed sequence-retrieval programming framework.
Using Workbench tool in the package, an input FASTA database is transformed into an index database containing a series of binary indexes (shown in Fig. 1a): two types of entry index (protein entry and peptide entry) and two types of mass index (peptide mass and ion mass). The former two types help to quickly locate the specific position to read fixed segments in the peptide and protein sequence.

The latter two types, typically designed for database searching and modification discovering (Tang et al., 2005), index a protein database by two types of sorted masses: the calculated no-modification masses of the peptides obtained from digestion of the proteins and the masses of theoretical no-modification MS/MS fragment ions such as b-ions of each of the peptide. The detailed data structure and generating process of the index are provided on the web pages named ‘Format’ and ‘Q&A’.

Once the protein database has been fully indexed, searching database and scoring an MS/MS spectrum is extremely rapid. To illustrate the improvement, the IPI RAT database, which is 21 MB containing 33,379 protein entries, is indexed by IndexToolkit. As we know, a frequent but time-consuming step for database searching is to find candidates whose mass fall within a mass error tolerance of the specified mass value. We evaluate the performance of using IndexToolkit by comparing the time running on un-indexed and indexed database (Fig. 1b). Figure 1b shows that searching with the indexes of IndexToolkit is many times faster than directly searching the raw FASTA database.

3 DISCUSSION

The CPU and memory usage may be high when indexing, but once indexing is completed, the future queries will usually use much less system resources to achieve a much higher performance. Index is a proper data structure to swiftly find all results satisfied specified criteria. However, it is not suitable for all cases. Because the cost to update indexes may be higher than improvements of querying performance, it is not recommended to index if the database changes frequently. In addition, large databases require large amount of disk space for storing their indexes. For IndexToolkit, the peptide-mass indexes may use ~3 times the disk space as the original database, and the ion-mass indexes may use ~30 times as the disk space. Nevertheless, we still suggest databases searching software run on the indexed database for high-throughput protein identification, as disk space is very cheap.

In sum, IndexToolkit can improve greatly the speed of protein database search engines, and is recommended for high-throughput protein identification. We have applied IndexToolkit to pFind (Li et al., 2005), which is a recently developed search engine with a novel scoring function for peptide identification. The new version of pFind integrated with IndexToolkit is about five times faster than the version without indexing. In our future work we will integrate IndexToolkit into other open-source search engines such as XTandem, and include more supporting database formats and additional tools for supporting protein identification.

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