Multiple Sequence Alignment using Anytime A*

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Alignment of multiple DNA or protein sequences is a central problem in computational biology. To create an alignment, gaps are inserted into sequences to shift characters to matching positions and a scoring function is used to rank the biological plausibility of alignments. Multiple sequence alignments are used to identify homologies among different species that reveal evolutionary history from a common ancestor. They are also used to discover genetic causes of certain diseases and to predict protein structure, which has significant importance in the design of drugs.

The multiple sequence alignment problem can be formalized as a shortest-path problem through a d-dimensional lattice, where d is the number of sequences to be aligned (Gusfield 1997). Dynamic programming is the traditional approach to constructing optimal alignments. Improved performance has recently been achieved using A*. However, the multiple alignment problem presents a difficulty for the classic A* algorithm. Its branching factor of the number of sequences d is so large that the size of the open list dramatically exceeds the number of nodes A* must expand to find an optimal solution.

Two solutions to this problem have been proposed in the literature. Yoshizumi et al. (2000) describe an extension of A*, called A* with Partial Expansion (PEA*). Instead of generating all successors of a node when it is expanded, PEA* inserts only the most promising successors into the open list. The “partially expanded” node is re-inserted into the open list with a revised f-cost equal to the least f-cost of its unexpanded successors, so that it can be re-expanded later. Use of this technique dramatically reduces the size of the open list, and PEA* can solve larger multiple sequence alignment problems than A*.

Unfortunately, the reduced space complexity of PEA* is achieved at the cost of node re-expansion overhead. The tradeoff between space and time complexity is adjusted by setting a “cutoff value” C, which determines which successor nodes to add to the open list.

Another way to reduce the size of the open list is to prune nodes from the open list if their f-cost is equal to or greater than a previously established upper bound, since such nodes will never be expanded by A*. This approach was first proposed by Ikeda and Imai (1999), who called it enhanced A* (EA*). One way to obtain an upper bound is to use the solution found by weighted A* search using a weight w > 1 in the node evaluation function f(n) = g(n) + wh(n). Ikeda and Imai suggested this method of obtaining an upper bound, but did not report experimental results for it.

In this abstract, we describe a third approach to reducing the size of the open list. In this approach, we also use weighted A* search to quickly find a solution that provides an upper bound that can be used to prune the open list. But because the first solution found may not be optimal, we continue the weighted search in order to find a sequence of improved solutions that eventually converges to an optimal solution. This also provides a sequence of improved upper bounds that can further prune the open list. We call this strategy Anytime A* (Hansen & Zilberstein 1996). Anytime A* refines both an upper bound, corresponding to the cost of the best solution found so far, and a lower bound, given by the unexpanded node with the least unweighted f-cost. Both bounds approach each other until convergence to a provably optimal solution. Before convergence, the difference between the two bounds gives an error bound on the quality of the currently available solution. Pseudocode for the algorithm is given at the end of the paper. The open list is pruned in lines 10 through 12.

Figures 1 and 2 compare the performance of Anytime A* (ATA*) to A* with Partial Expansion and Enhanced A* (where Enhanced A* uses the first solution found by weighted A* as an upper bound to prune the open list). The PAM250 cost matrix is used with a gap cost of 8. All three algorithms require dramatically less memory than conventional A* in solving the multiple sequence alignment problem, allowing a larger number of sequences to be aligned.

Figure 1 compares their performance in aligning eight sequences from a highly similar set of sequences used in earlier experiments (Ikeda & Imai 1999; Yoshizumi, Miura, & Ishida 2000). On average, Anytime A* runs more than 7 times faster and stores only 26% more nodes than PEA* using a cutoff of C = 0. When PEA* uses a cutoff of C = 50, it stores 40% more nodes than Anytime A* and still runs 18% slower on average. Enhanced A* performs best on this test set. It runs 20% faster than Anytime A* and stores 4% fewer nodes.

Figure 3 compares the performance of the algorithms in aligning five sequences from a set of dissimilar sequences used in earlier experiments (Kobayashi & Imai 1998). For
A* is optimal or close to it—a best-case scenario for En-
n this test set. Anytime A* is only 2.4% slower than enhanced
A* and stores 40% fewer nodes. Its better performance on
the second test set is explained as follows. Because the se-
quences in the first test set are very similar and the heuristic
is extremely accurate, the first solution found by weighted
A* and stores 40% fewer nodes. Its better performance on
this test set, Anytime A* is only 2.4% slower than enhanced
A* and stores 40% fewer nodes. Its better performance on
the second test set is explained as follows. Because the se-
quences in the first test set are very similar and the heuristic
is extremely accurate, the first solution found by weighted
A* is optimal or close to it—a best-case scenario for En-
hanced A*. Because the sequences in the second set are
dissimilar and the heuristic is less accurate, the first solu-
tion found by weighted A* is usually not optimal. Anytime
A* continues to find better solutions that improve the upper
bound, which in turn improves memory-efficiency. Figure 2
illustrates its anytime behavior by showing how the upper
and lower bounds gradually converge.

Our experimental results show that the sequence of im-
proved solutions found by Anytime A* provide a dynamic
upper bound that keeps its memory requirements close to
the minimum number of nodes that must be expanded to find
an optimal solution. Anytime A* is also more memory-efficient
than PEA* unless the latter uses the most aggressive cutoff
of $C = 0$, in which case the node re-expansion overhead
of PEA* slows it considerably. Anytime A* is also more
memory-efficient than Enhanced A* when aligning dissimi-
lar sequences. Anytime A* has an additional advantage over
both algorithms. Because it finds a sub-optimal alignment
quickly and continues to improve the alignment with addi-
tional computation time, it offers a tradeoff between solution
quality and computation time that can prove useful when
finding an optimal alignment is infeasible.

**Pseudocode of Anytime A**

```plaintext
1 $g(s) \leftarrow 0$, $f(s) \leftarrow g(s) + w \times h(s)$
2 OPEN $\leftarrow \{s\}$, CLOSED $\leftarrow \emptyset$, bound $\leftarrow \infty$
3 while OPEN $\neq \emptyset$
4   $n \leftarrow \text{arg min}_{x \in \text{OPEN}} \{f(x) \mid x \in \text{OPEN}\}$
5   OPEN $\leftarrow$ OPEN $\setminus \{n\}$
6   CLOSED $\leftarrow$ CLOSED $\cup \{n\}$
7 if $n$ is a goal node then
8     bound $\leftarrow g(n) + h(n)$
9     Output solution and bound
10 for each $x \in \text{OPEN and}$
11   $g(x) + h(x) \geq \text{bound}$ do
12     OPEN $\leftarrow$ OPEN $\setminus \{x\}$
13 for each $n_i \in \{x \mid x \in \text{Successors}(n),$
14   $g(n) + c(n, x) + h(x) < \text{bound}\}$ do
15   if $n_i \notin \text{OPEN} \cup \text{CLOSED}$ or
16     $g(n_i) > g(n) + c(n, n_i)$ then
17     $g(n_i) \leftarrow g(n) + c(n, n_i)$
18     $f(n_i) \leftarrow g(n_i) + w \times h(n_i)$
19     OPEN $\leftarrow$ OPEN $\cup \{n_i\}$
20 if $n_i \in \text{CLOSED}$ then
21     CLOSED $\leftarrow$ CLOSED $\setminus \{n_i\}$
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**References**

Gusfield, D. 1997. *Algorithms on Strings, Trees, and Se-
quen: Computer Science and Computational Biology*. Cambridge University Press.


Ikeda, T., and Imai, H. 1999. Enhanced A* algorithms for multiple alignments: optimal alignments for several se-
