Optimal Detection of Sequence Similarity by Local Alignment

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The statistical properties of local alignment algorithms with gaps are analyzed theoretically for uncorrelated and correlated random sequences. In the vicinity of the log-linear phase transition, the statistics of alignment with gaps is shown to be characteristically different from that of gapless alignment. The optimal scores obtained for uncorrelated sequences obey certain robust scaling laws. Deviation from these scaling laws signals sequence homology, and can be used to guide the empirical selection of scoring parameters for the optimal detection of sequence similarities. This can be accomplished in a computationally efficient way by using a novel approach focusing on the score profiles. Furthermore, by assuming a few gross features characterizing the statistics of underlying sequence-sequence correlations, quantitative criteria are obtained for the choice of optimal scoring parameters: Optimal similarity detection is most likely to occur in a region close to the log side of the loglinear phase transition.

Keywords: sequence alignment; homology; optimization; phase transition

1 INTRODUCTION

Sequence alignment is a vital tool in molecular biology. It has been used extensively in discovering and understanding the functional and evolutionary relationships among genes and proteins [10, 37]. There are two basic types of alignment algorithms: algorithms without gaps, e.g., BLAST and FASTA [1], and algorithms with gaps, e.g., the Needleman-Wunsch algorithm [27] and the Smith-Waterman algorithm [30]. Gapless alignment is widely used in large-scale database searches because the algorithms are fast [1], the results depend only weakly on the choice of scoring systems [2], and the statistical significance of the results is well-characterized [4, 21, 22]. However, gapless alignment is not sufficient for the detection of *weak* sequence similarities [29]. For the detailed analysis of such sequences, algorithms with gaps are necessary [13, 35, 37]. Advancing our understanding of the

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statistics of gapped alignment could therefore be critical to the wider usage of these more powerful tools.

A notorious difficulty for any alignment is the selection of scoring schemes and/or parameters: In a generic sequence matching problem, a score is assigned to each alignment of given sequences, based on the total number of matches, mismatches, gaps, etc. Maximization of this score defines an optimal alignment. However, it is well known that the optimal alignment of given sequences strongly depends on the particular scoring scheme and/or parameters used. Consequently, the *fidelity* of an alignment, i.e., the extent to which the alignment captures mutual correlations among the aligned sequences, can depend strongly on the choice of scoring parameters. Understanding the influence of these parameters on the resulting alignment and choosing the appropriate parameters are therefore important for the proper application of these algorithms. This requires the knowledge of the statistics of alignment results, which has been obtained only for gapless alignments [2, 4, 21, 22]. For alignments with gaps, appropriate parameters have so far been chosen mostly by trial and error, although there have been systematic efforts to establish a more solid empirical footing [6, 32]

Recently [18, 11] we have analyzed the statistical properties of *global* alignment with gaps. Such algorithms align sequences of similar lengths in their entirety. By exploiting mathematical analogies to certain well-studied problems of statistical mechanics [19, 16, 17, 23], we have obtained a quantitative description of the global alignment statistics for mutually *uncorrelated* as well as *correlated* sequence pairs. Here we extend the analysis to *local* alignment algorithms [30] which find the best match between contiquous subsequences, subject to (finite) penalties for gaps and mismatches. For uncorrelated random sequences, i.e., for independent sequences with iid or Markov letters, it is well known that depending on the choice of scoring parameters, the length of the optimal subsequence alignment depends either linearly or logarithmically on the total length of the sequences [34, 5]. A phase transition line separates the space of scoring parameters into two regimes: the "linear phase" for small gap and mismatch costs, and the "log phase" for large penalty costs. It is clear that local alignment deep in the linear phase is equivalent to global alignment (and hence described by the results of our previous studies). On the other hand, the log phase at high gap penalty becomes indistinguishable from the log phase of gapless alignment. Indeed there have been extensive empirical efforts [31, 7, 25, 39, 40, 3] to characterize the statistics of the log phase of gapped algorithms by an effective description

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as gapless alignment with modified parameters. While this approach is reasonable far away from the phase transition, it becomes questionable as the phase transition line is approached, since the linear phase itself is completely different from the log phase. On the other hand, the loci of scoring parameters for optimal similarity detection appear to lie in the log phase close to the phase transition line, according to recent empirical studies by Vingron and Waterman [32]. Hence, understanding the log-linear phase transition is crucial for optimizing the detection of sequence similarity and quantifying its statistical significance.

In this work, we apply the well-established theory of phase transition [9] to the log-linear transition of gapped local alignment. We find various statistical properties at and in the vicinity of the transition to be governed by scaling laws analogous to those recently discovered for global alignment [18, 11]. The transition turns out to differ qualitatively and quantitatively from its counterpart in gapless alignments. Our results lead to quantitative criteria for the optimal choice of scoring parameters, given certain gross statistical characteristics of the expected sequence correlation. In particular, they explain why optimal parameters for weakly correlated sequence pairs are in the vicinity of the phase transition line as observed by Vingron and Waterman [32]. Also emerging from this work is a versatile method to detect sequence correlation and characterize its statistical significance empirically for sequences with a priori unknown correlations.

2 REVIEW OF ALIGNMENT ALGORITHMS

We study the Smith-Waterman (SW) local alignment algorithm applied to a pair of long nucleotide sequences \mathcal{P}_1 and \mathcal{P}_2 , with lengths $N_1 \simeq N_2 \gg 1$. Let $P_{n,i} \in \{A, T, G, C\}$ be the i^{th} element of the sequence \mathcal{P}_n . A particular alignment consists of an ordered set of pairings of two elements $(P_{1,i}, P_{2,j})$, or of an element with a gap, for any contiguous subsequence of length $\ell_1 \leq N_1$ in sequence \mathcal{P}_1 and length $\ell_2 \leq N_2$ in sequence \mathcal{P}_2 [see Fig. 1]. The simplest scoring system assigns a positive score of +1 if the two elements paired together are identical, and a negative score of $-\mu$ if the two are different. Each pairing of an element with a gap is penalized with a negative score $-\delta$. (In the simple case considered here, we shall not distinguish between gap initiation and gap extension. It is then sufficient to consider only the region $2\delta \ge \mu^1$.) The sum of the scores of all individual pairings of a given alignment is the total score for that alignment. An optimal alignment is one for which the total score is maximized for a given set of scoring parameters (μ, δ) . The SW algorithm uses the dynamic programming method to find the optimal alignment of all possible subsequences. Key to the algorithm is the unique representation of an alignment by a directed path (see Ref. [27] and Fig. 1). Let us briefly recall the algorithm below, cast in a slightly different notation to facilitate the subsequent analysis.

Consider the alignment lattice shown in Fig. 1. We label each lattice point by the coordinate (x, z), with the lower tip of the lattice anchored at (0, 0). The highest total score of all alignment paths ending at a point (x, z) is denoted by h(x, z). Given h(x, z) and h(x, z - 1) for all x, h(x, z + 1)

$$\{P_I\}$$
 G G — A C A T A

$$\{P_2\}$$
 C **G T A T A** G

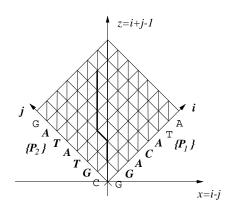


Figure 1: One possible local alignment of the two nucleotide sequences, $\mathcal{P}_1 = GGACATA...$ and $\mathcal{P}_2 = CGTATAG...$ The aligned subsequences are shown in boldface, with 4 pairings (three matches, one mismatch) and one gap. This alignment can be represented uniquely as a directed path on the alignment lattice; each left (right) turn of the path correspond to a gap insertion in sequence \mathcal{P}_1 (\mathcal{P}_2).

can be computed from the recursion relation

$$h(x, z+1) = \max \left\{ \begin{array}{c} h(x+1, z) - \delta \\ h(x-1, z) - \delta \\ h(x, z-1) + u(x, z) \\ h_0 \end{array} \right\}, \quad (1)$$

with δ being the gap insertion cost, u being the match/mismatch score to be specified below, and h_0 being a cutoff score. The SW algorithm has $h_0 = 0$, which effectively deletes the segment of the alignment path connecting (0,0) and (x,z) if $h(x,z) \leq 0$. In contrast, the global alignment algorithm of Needleman and Wunsch has no cutoff, corresponding to the limit $h_0 \to -\infty$. Also, gapless local alignment (with $\delta \to \infty$) corresponds to the limit of the recursion relation (1) involving only one value of x, say x = 0. The scoring function u(x, z) gives the match/mismatch score of aligning the elements $P_{1,i}$ with $P_{2,j}$, with x = i - j and z = i + j - 1. For simplicity, we use in this study the form

$$u(x=i-j, z=i+j-1) = \begin{cases} 1 & \text{if } P_{1,i} = P_{2,j} \\ -\mu & \text{if } P_{1,i} \neq P_{2,j} \end{cases} . (2)$$

More elaborate forms of the scoring function are easily incorporated and do not change key results of this study.

If z and x are regarded as "time" and "space" variables, respectively, the recursion relation (1) can be viewed as a "dynamical process" describing the time evolution of the one-dimensional "score profile" h(x, z). (Similarly, gapless local alignment involving only the site x = 0 corresponds to the "zero-dimensional" limit.) This dynamic analogy will be pivotal in guiding the ensuing analysis.

3 GLOBAL ALIGNMENT

¹For $2\delta < \mu$, it is always favorable to replace a mismatch by two gaps, so the outcome of an alignment becomes independent of μ .

3.1 Statistics of Uncorrelated Sequences: Universal Scaling Laws

Let us first review some of the relevant results we previously obtained for the global alignment of random sequence pairs [18]. In Fig. 2(a), we show several representative constantz slices of the score profile h(x) obtained by iterating Eq. (1) with $(\mu, \delta) = (0.5, 2.0)$. The alignment algorithm is applied to one pair of random sequences each of length N = 10000. Results are shown for a central rectangular region² of the alignment lattice, $-X/2 \leq x \leq X/2$ and $X/2 \leq z \leq 2N - X/2$ with $X \leq N$, starting from the initial condition h(x, z = X/2) = 0. It will be convenient to use a shifted time-like variable, $t \equiv z - X/2$. In Fig. 2(a), we see a series of disorderly score profiles, with the "spatial" average

$$h(t) = X^{-1} \sum_{x=-X/2}^{X/2} h(x,t)$$
(3)

advancing steadily in t. For large t, we obtain the linear dependence, $h(t) = v_0(\mu, \delta)t$ (not shown). The value of the rate v_0 itself is not important for global alignment. (Thus, h(t) could as well be *decreasing* linearly in t.) More significant is the spatial variation in the profile which always increases for increasing t. This is more clearly seen by plotting the effective width of the profile, w(t), defined as

$$w^{2}(t) = \frac{1}{X} \sum_{x=-X/2}^{X/2} \left[h(x,t) - h(t)\right]^{2}, \qquad (4)$$

or alternatively, the difference between the highest and lowest point of the profile, $\Delta h(t) = h_{\max}(t) - h_{\min}(t)$; see Fig. 2(b).

The roughness of the profile, as quantified by either w(t)or $\Delta h(t)$, is an important characteristic of the alignment. since it indicates how strongly the optimal alignment dominates over the suboptimal alignments. The score profiles of Fig. 2(a) show the weak dominance of the optimal alignment and the existence of a large number of suboptimal alignments. The statistics of these suboptimal alignments has been recognized recently as a valuable tool in sequence alignment; for an interesting recent exposition, see Ref. [33]. From Fig. 2(b), it appears that the roughness grows with a sub-linear power in t. This is verified in Fig. 3(a), where we show the ensemble average $\overline{w}(t)$ for different sets of scoring parameters (μ, δ) , each curve averaged over 1000 pairs of uncorrelated random sequences. (Throughout the text, we use overbars to denote averages of an ensemble of random sequence pairs.) It is seen that for large t, the width obeys the asymptotic scaling law

$$\overline{w}(t) = B(\mu, \delta) t^{\omega}.$$
(5)

Different parameter choices only affect the prefactor B, but not the exponent $\omega \approx 1/3$. The same scaling law (with a larger coefficient) is found also for $\overline{\Delta h(t)}$.

The scaling law (5) is an example of the "universal scaling phenomena" well studied in statistical mechanics [9].

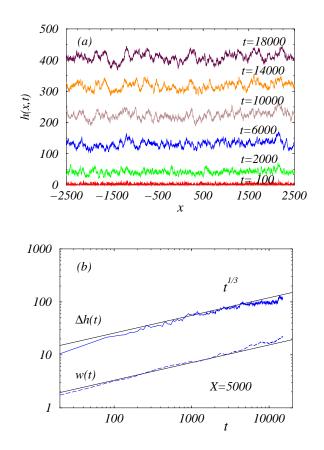


Figure 2: (a) A typical series of score profiles h(x,t) obtained from the global alignment of a pair of uncorrelated random sequences; parameters used are $(\mu, \delta) = (0.5, 2.0)$. (b) Gradual increase in the "roughness" of the profile, as characterized by either the width w(t) or the range $\Delta h(t)$ over the range of x shown in (a). The straight lines indicate the suggested power law dependence.

The exponent ω is a very robust characteristics of random sequence alignment. It quantifies the roughness of the profile and hence the dominance of the optimal alignment. It does not depend on details of the scoring function, (e.g., whether or not gap initiation and extension are differentiated) but only on a few qualitative characteristics such as the number of sequences aligned or the type of the correlations between them.

In Ref. [18], we have given arguments suggesting that the result $\omega = 1/3$ is exact for the global alignment of random sequence pairs. Our approach is based on the close analogy of the recursion relation (1) and discrete models of kinetic surface growth studied extensively in statistical mechanics in the past decade³. In these growth models, h(x, t) describes the height profile of a one-dimensional surface at time t. The growth is driven by a stochastic process that governs the deposition and removal of material on the surface and is described by the random variable $\eta(x, t) = \frac{1}{2}u(x, t) - \delta$. The large scale behavior of the profile h(x, t) generated by the growth model is well described by the *differential* growth

²Due to boundary effects arising from the diamond-shaped alignment lattice (Fig. 1), the total score h(x,t) certainly decreases (quadratically on average) as one moves far away from the center at x = 0. To remove these spurious effects, we focus our attention only to the central strip of the alignment lattice, e.g., for $-X/2 \le x \le X/2$ and $X/2 \le z \le 2N - X/2$, with $X \le N$. All results reported here are obtained using this strip geometry. For long sequences, the statistics of the score profile obtained from the strip is indistinguishable from that obtained with the full alignment lattice.

 $^{^{3}}$ General reviews of the kinetic growth problem and its relation to the problem of first passage percolation can be found in Refs. [24] and [15].

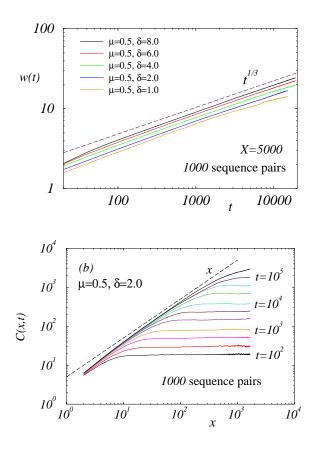


Figure 3: (a) The ensemble averaged roughness (width) for different sets of scoring parameters, each averaged over 1000 pairs of randomly generated sequences of length N = 10,000. The dashed straight line indicates the expected power law form. (b) The ensemble averaged spatial correlation function C(x,t) for $(\mu, \delta) = (0.5, 2.0)$. The dashed line indicates the expected scaling behavior at large t. For small t's, C(x,t) saturates to the order of $\overline{w}^2(t)$.

equation [20]

$$\frac{\partial h}{\partial t} = \nu \frac{\partial^2 h}{\partial x^2} + \frac{\lambda}{2} \left(\frac{\partial h}{\partial x}\right)^2 + \eta(x, t), \tag{6}$$

with the coefficients ν and λ being functions of δ and μ . Eq. (6) is known as the Kardar-Parisi-Zhang equation, closely related to the noise-driven Burgers' equation [12]. If $\eta(x,t)$ is an uncorrelated Gaussian random variable, the stationary state $(t \to \infty)$ of the surface can be obtained exactly [24], resulting in the Gaussian equal-time distribution

$$P\{h(x,t\to\infty)\} \propto e^{-\frac{1}{2D}\sum_{x}[h(x+1,t)-h(x,t)]^2},$$
 (7)

and a coefficient $D(\mu, \delta)$. Approaching the steady state, one has $\overline{w}(t \gg 1) = B t^{\omega}$, with $B \approx D^{2/3}$ and $\omega = 1/3$ exactly. For sequence alignment, the variables $\eta(x, t)$ and $\eta(x', t')$ at different points are not independent given their definition above. One can verify for example that the higher moments of η are long-range correlated. However, as argued in Ref. [18] and [Hwa and Lässig (to be published)], these correlations do not affect the asymptotic scaling behavior⁴. The close correspondence of the numerical results on w(t) (Fig. 3(a)) and the exact result $\omega = 1/3$ supports this conclusion. An independent check is to measure directly the equal-time correlation of h, $C(x,t) = \overline{[h(x + x',t) - h(x',t)]^2}$. From Eq. (7), we expect to have C(x,t) = D |x| for sufficiently large t (or sufficiently small x). For finite t, C(x,t) must eventually saturate, to a value $\sim \overline{w}^2(t)$ for $x > O\left(t^{2/3}\right)$.

The numerically obtained forms of C(x, t) are shown in Fig. 3(b) for $(\mu, \delta) = (0.5, 2.0)$. The results are in good agreement with the anticipated form. Similar results are obtained for other values of the scoring parameters.

3.2 Correlated Sequences: Similarity Detection from the Score Profile

To illustrate how the above knowledge can be used for the purpose of similarity detection, we next describe the score profile for the global alignment of *correlated* sequences. Sequence correlations are obtained by first generating a random "ancestor sequence", and then making imperfect replicas \mathcal{P}_1 and \mathcal{P}_2 . The degree of sequence-sequence correlation is quantified by the average number of point substitutions per base p_s and the average number of indels per base p_t made in the replication of each daughter sequence. (See Refs. [18, 11] for details.) In Fig. 4(a), we show the score profile h(x,t) and the range $\Delta h(t)$ for a pair of heavily mutated daughter sequences with $p_s = 40\%$ and $p_t = 20\%$, corresponding to a fraction of < 30% conserved elements. The score profiles obtained are very different from the profiles characteristic of uncorrelated random sequences shown in Fig. 2(a). For correlated sequences (Fig. 4(a)), a peak emerges from the disordered background after some time. The height of the peak advances steadily at a rate vwhich exceeds the growth rate of the background v_0 . Thus, the peak gradually broadens to engulf the entire profile. The dominance of the central peak reflects the existence of strongly preferred alignments for correlated sequences, in marked contrast to the alignment of random sequences⁵.

The existence of the peak can be taken as a manifestation of inter-sequence correlations. The strength of sequence correlation detected is quantified by the difference between the growth rate of the peak and the background, $\varepsilon \equiv v - v_0$. The magnitude of ε is clearly dependent on the mutation rates p_s and p_t , but also depends on the scoring parameters μ and δ . The functional form of $\varepsilon(\mu, \delta; p_s, p_t)$ has recently been investigated in detail [11] and will not be addressed here.

A peak in the score profile is discernible from the background only if the height of the peak, of the order $\varepsilon \cdot t$ after tsteps, exceeds the roughness of the background $\sim B \cdot t^{1/3}$ (see Fig. 4(b)). Hence, using the score profile approach, one can detect correlations between sequences if their lengths exceed the threshold length

$$t_c(\mu, \delta; p_s, p_t) \sim \left[B(\mu, \delta) / \varepsilon(\mu, \delta; p_s, p_t) \right]^{3/2}.$$
 (8)

Minimization of t_c (for a given sequence pair) is a natural empirical criterion for optimal similarity detection, yielding a preferred choice of scoring parameters $\delta^*(\mu; p_s, p_t)$. A

 $^{^4{\}rm More}$ detailed discussions are given in Ref. [8] in the context of a number of closely related physics problem.

⁵In statistical mechanics, one considers the free energy, which is the negative of the score. The score profile discussed here is a measure of the "energy landscape". A peak in the score profile corresponds to a valley or a "funnel" in the energy landscape. It has been suggested that the energy landscape of heteropolymeric systems such as a protein has the funnel shape [28]. The score profile (Fig. 4(a)) obtained here is the first known example of a large heteropolymeric system for which the suggested landscape is directly observed.

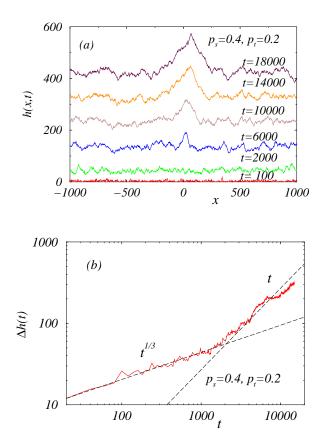


Figure 4: (a) The score profiles obtained from the global alignment of a pair of weakly correlated sequences (with conserved fraction < 30%) are characteristically different in appearance from those of the uncorrelated sequences shown in Fig. 2(a). (b) The roughness corresponding to the profiles of (a) begins to deviate from the $t^{1/3}$ behavior at $t \simeq 2000$.

more fundamental criterion for choosing the optimal scoring parameters is to maximize the *fidelity* of the alignment, i.e., the extent to which the optimal alignment reconstructs the ancestor sequence [18]. The dependence of fidelity on scoring parameters has been studied in detail in Ref. [11]. It was found that maximizing the fidelity is closely related (and equivalent for practical purposes) to minimizing the threshold length t_c . Thus, the empirical criterion based on the score profile is indeed a versatile way of locating the optimal parameters.

The above strategy of similarity detection is close in spirit to the well-known "shuffling method", which compares the alignment score of a given sequence pair to the distribution of scores obtained from aligning the ensemble of randomly shuffled sequences (for the same set of scoring parameters). However, by making use of the spatiallyextended score profile and its time evolution (as opposed to keeping track of only the optimal score), we have demonstrated that this comparison can be accomplished by one single alignment. This is a drastic improvement over the shuffling method, which requires the generation and alignment of an ensemble of sequences. It should thus be possible to minimize $t_c(\mu, \delta)$ empirically for sequence pairs with a priori unknown correlation, since the value of t_c can be obtained directly from the onset of score peak shown in Fig. 4(b), without any assumption on the nature of sequence correlations. It would be particularly interesting to combine this approach with the efficient ensemble and parametric alignment algorithms [36, 14, 38] which find optimal scores for all parameters.

4 LOCAL ALIGNMENT WITH GAPS

We now describe the statistical properties of local alignment, with $h_0 = 0$ in the recursion relation (1). Local alignment is necessary since often only a subsequence of one sequence is correlated with a subsequence of another. Let the lengths of the correlated subsequences be $\ell_1 \approx \ell_2 \approx \ell$. If the positions of the subsequences were known, the correlations would be detectable by global alignment of these subsequences if $\epsilon \ell > B \ell^{1/3}$, i.e., if $\ell > t_c$. However, if global alignment is applied to the entire sequences, the background noise $BN^{1/3}$ will in general overwhelm the score signal of the correlations $\epsilon \ell$. The advantage of local alignment is that by cutting off the length of the aligned segment, it limits the background roughness to a *finite* value such that the correlation peak can still be detected.

4.1 Uncorrelated Sequences: the Log-Linear Phase Transition

We first discuss the background, i.e., the local alignment of uncorrelated random sequence pairs. Unlike global alignment, the outcome of local alignment depends critically on the value of v_0 , i.e., the growth rate of the average score of the corresponding global alignment problem. For scoring parameters (μ, δ) in the regime where $v_0(\mu, \delta) > 0$, the score h(x,t) grows without bound also for local alignment, and the existence of a cutoff score $h_0 = 0$ is immaterial in the limit of large t. Thus the asymptotic properties of local and global alignments are identical in this regime. (It acquires the name "linear phase" since the average score $\overline{h}(t)$ advances linearly by definition.) For $v_0(\mu, \delta) < 0$, on the other hand, the average $\overline{h}(t)$ saturates quickly to a constant value, and the largest value $h(x, t' \leq t)$ scales as $\log(t)$ due to exponentially rare events. As has been pointed out by Arratia and Waterman [unpublished], the condition⁶

$$v_0(\mu,\delta) = 0 \tag{9}$$

defines the phase transition line separating the two regimes (see Fig. 5). The statistical properties in the vicinity of this phase transition have recently been studied in the context of some related physics problems [26]. We will briefly summarize the results below.

The phase transition is very similar to that of gapless local alignment, except that the $t^{1/2}$ score dependence at the transition is replaced by

$$h_c(t) \equiv \overline{h}(t)_{\mu=\mu_c} = b(\delta) t^{1/3}, \qquad (10)$$

along the critical line $\mu_c(\delta)$ for large t, with a coefficient $b(\delta) \sim B(\mu_c(\delta), \delta)$ dependent on the location along the phase transition line. Slightly away from the critical line on the log side, i.e., for $\tilde{\mu} \equiv \mu - \mu_c \gtrsim 0$ where $v_0 \propto -\tilde{\mu} \lesssim 0$, $\overline{h}(t)$

⁶However, finite-size effects [9] may cause a slight shift in the apparent transition point. For example, the fits of Figs. 6 and 7 found $\mu_c \approx 0.7085$ for $\delta = 2.0$, whereas $v_0 = 0$ occurred at $\mu \approx 0.7040$ for $\delta = 2.0$.

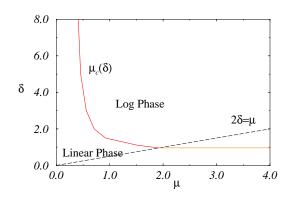


Figure 5: The critical line $\mu_c(\delta)$ separating the log and the linear phases is obtained from the condition $v_0(\mu_c, \delta) = 0$. (The critical line becomes independent of μ below the line $2\delta = \mu$.)

is indistinguishable from $h_c(t)$ for small t, saturating eventually to a constant value $h_{\text{sat}} \approx b^{3/2}/|v_0|^{1/2}$. The saturation scale t_{\times} is obtained from the condition $h_c(t = t_{\times}) \sim h_{\text{sat}}$, yielding

$$t_{\times} \sim (b/|v_0|)^{3/2} \propto |\widetilde{\mu}|^{-3/2}.$$
 (11)

This is the length of the the optimally aligned subsequences selected by local alignment. For sequence lengths N much longer than t_{\times} , it is reasonable to approximate the score statistics by that of gapless local alignment, as has been attempted previously [31, 7, 25, 39, 40, 3]. However, because t_{\times} diverges as the critical point ($\tilde{\mu} = 0$) is approached, the gapless approximation becomes invalid for $t_{\times} > N$, or for $\tilde{\mu} < N^{-2/3}$. On the other side of the critical line where $\tilde{\mu} < 0$ and $v_0 > 0$, $\bar{h}(t)$ again equals $h_c(t)$ for $t \lesssim t_{\times}$, before changing to the linear dependence for larger values of t. The statistics of global alignment applies to the linear phase on scales $t > t_{\times}$.

The above behavior of $\overline{h}(t)$ can be summarized by the homogeneous scaling relation

$$\overline{h}(t;\tilde{\mu}) = |\tilde{\mu}|^{-1/2} f_{\pm}(t\,|\tilde{\mu}|^{3/2}),\tag{12}$$

with the two branches of the scaling functions $f_{\pm}(x)$ having the limiting forms $f_{\pm}(x \ll 1) = b x^{1/3}$, $f_{+}(x \gg 1) \rightarrow \text{const}$ for $\tilde{\mu} > 0$ and $f_{-}(x \gg 1) \propto x$ for $\tilde{\mu} < 0$ ⁷. Such scaling relations are widely encountered in physical systems undergoing continuous phase transitions and have been studied extensively by the modern theory of critical phenomena [9].

In Fig. 6(a), we show the score average $\overline{h}(t)$ for local alignment of 1000 random sequence pairs of length 10000, for $\delta = 2$ and various μ 's taken $\pm 5\%$ around the critical value $\mu_c \approx 0.7085$. The anticipated scaling form (12) can be checked by multiplying the horizontal axis of Fig. 6(a) by a factor $|\tilde{\mu}|^{3/2}$ and the vertical axis by $|\tilde{\mu}|^{1/2}$ individually for each curve. The result is shown in Fig. 6(b). The 11 curves displayed in Fig. 6(a) (each containing 10⁵ data points) collapse into two branches, corresponding to the two branches of the scaling function f_{\pm} . The upper branch (f_{-}) describes

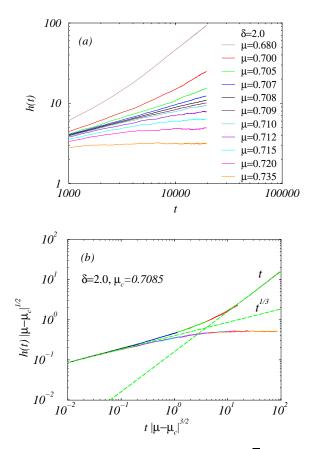


Figure 6: (a) The ensemble averaged score $\overline{h}(t)$ obtained from the local alignment of random sequence pairs, with $\delta = 2.0$, and various values of μ slightly above and below the critical value $\mu_c(\delta = 2.0) \approx 0.7085$. (The order of the curves corresponds to the order shown in the legend, with the top curve having the smallest value of μ and the bottom curve having the largest value of μ .) Each curve is averaged over 1000 pairs of sequences of length 10,000 each. (b) The curves in (a) plotted according to the homogeneous scaling form (see text). The dashed lines indicate the anticipated power law forms in the two different regimes.

the crossover of the critical behavior $h_c(t)$ to the linear behavior, and the lower branch (f_+) describes the crossover of the critical behavior towards saturation. The only fitting parameter for this data collapse is the location of the transition point μ_c .

A similar scaling relation exists for the roughness of the score profile. In Fig. 7(a), the width $\overline{w}(t)$ is plotted for various values of μ , approaching the phase transition from the log side. The data can be collapsed by the same transformation as in Fig. 6(b), with the same fitting parameter μ_c . The results (Fig. 7(b)) show clearly that $w(t; \tilde{\mu})$ has the same form as $\overline{h}(t; \tilde{\mu})$ in the log phase, i.e., the lower branch of Fig. 6(b). In particular, $w(t > t_{\times})$ saturates to a constant value w_{sat} of the same order as h_{sat} , yielding

$$w_{\rm sat} \sim b t_{\times}^{1/3}.\tag{13}$$

This is just the expression (5) describing the score roughness of *global* alignment, evaluated at $t = t_{\times}$. It is a manifestation of the general result that local alignment corresponds to global alignment applied to the selected subsequences.

⁷ It should be noted that the scaling form (12) also describes the score $h_{\text{gapless}}(t;\tilde{\mu})$ of the zero-dimensional problem of gapless local alignment in the vicinity of its phase transition, but with modified exponents. The results $h_{\text{gapless}}(t;\tilde{\mu}) = |\tilde{\mu}|^{-1}g_{\pm}(t|\tilde{\mu}|^2)$, with $g_{\pm}(x \ll 1) \propto x^{1/2}, g_{\pm}(x \gg 1) \rightarrow \text{const}$ and $g_{-}(x \gg 1) \propto x$ can be straightforwardly verified; see Ref. [26] for a detailed discussion.

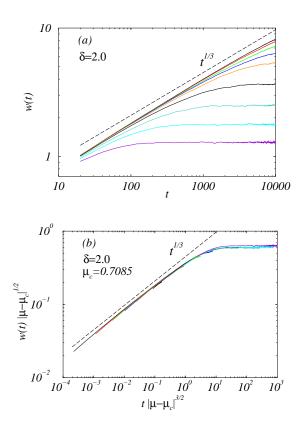


Figure 7: (a) The ensemble averaged roughness $\overline{w}(t)$ in the log phase, for $\delta = 2.0$ and $\mu = 0.709$, 0.710, 0.712, 0.715, 0.720, 0.735, 0.765, 0.825, 0.950 (from top to bottom). The sequences are the same as those used in Figs. 6. (b) The roughness curves plotted according to the scaling form.

4.2 Correlated Subsequences: Similarity Detection and Parameter Dependence

The existence of a phase transition in local alignment can be used to detect sequence-sequence correlations: Consider a pair of sequences \mathcal{P}_1 and \mathcal{P}_2 with mutually correlated subsequences located in the intervals from i_0 to $i_0 + \ell/2$ and from j_0 to $j_0 + \ell/2$, respectively. For concreteness, let the similarity of these subsequences be generated by the evolution mechanism described in Sec. 3.2, with the strength of inter-sequence correlations characterized by $\varepsilon(\mu, \delta; p_s, p_t)$. Our strategy for similarity detection is simply to choose the scoring parameters δ and μ such that $v_0(\mu, \delta) < 0$, i.e., the log phase is obtained if the sequences are uncorrelated, while keeping $v_0 + \varepsilon > 0$, so that the linear phase may be obtained instead for some duration of the correlated subsequences, $t_0 < t < \ell$. With this parameter choice, the profile h(x,t) will have a constant background roughness $w_{\rm sat}$ for $t < t_0 = i_0 + j_0 - 1$, and a score peak signaling subsequence correlations in the interval $t_0 < t < \ell$, once

$$(\epsilon + v_0) \cdot (t - t_0) > w_{\text{sat}}.$$
(14)

Optimal similarity detection is obtained by maximizing the peak-to-background ratio,

$$\sigma = (\epsilon + v_0)/w_{\rm sat} \tag{15}$$

which can be taken as a measure of the statistical significance of the correlations detected. Using the relations (13) and (11), we find $\sigma \sim (\varepsilon - |v_0|) \cdot |v_0|^{1/2}$, which is maximized at

$$v_0^* = v_0(\mu^*, \delta^*) = -\frac{\varepsilon}{2},$$
 (16)

with

$$\sigma^* \propto \varepsilon^{3/2} \sim t_c^{-1}. \tag{17}$$

Eqs. (16) and (17) are the central results of this study. Eq. (16) shows that for weakly correlated subsequences (i.e., $\varepsilon \to 0$), the optimal scoring parameters should be close to the log side of the phase boundary. This is a quantitative formulation of the empirical observation of Vingron and Waterman [32]. In addition to the choice of v_0 , Eq. (17) shows that t_c should still be minimized as discussed in Sec. 3.2 for global alignment. These two conditions uniquely determine the optimal parameters μ^* and δ^* . Using the optimal result, the condition (14) is reduced to $t-t_0 > t_c^* = t_c(\mu^*, \delta^*; p_s, p_t)$, which yields the minimal subsequence length ℓ for which correlations can be detected.

5 SUMMARY

In this study, we presented a statistical description of the Smith-Waterman local alignment algorithm, focusing on the properties near the log-linear phase transition line. We demonstrated how this knowledge can be exploited to provide quantitative criteria guiding the choices of alignment parameters for optimal detection of weak sequence correlations. The optimal values μ^* and δ^* are obtained in terms of the substitution rate p_s and the indel rate p_t characterizing the statistics of sequence correlations. By analyzing the evolution of the spatially-extended score profile, we are able to detect sequence correlations by a single run of the algorithm. This is a very efficient way to optimize the scoring parameters empirically, and may be useful in the alignment of a vast number of weakly correlated sequences.

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