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Exact Bounds for Interval and Fuzzy Functions Under Monotonicity Constraints, with Potential Applications to Biostratigraphy

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Abstract— The age of fossil species in samples recovered from a well that penetrates an undisturbed sequence of sedimentary rocks increases with depth. The results of biostratigraphic analysis of such a sequence consist of several age-depth values – both known with interval (or fuzzy) uncertainty – and we would like to find, for each possible depth, the interval of the possible values of the corresponding age. A similar problem of bounding an intervally (fuzzily) defined function under monotonicity constraint occurs in many other application areas. In this paper, we provide an efficient algorithm for solving this problem.

I. CASE STUDY: BIOSTRATIGRAPHY

A. Biostratigraphy is important

Biostratigraphy is concerned with the stratigraphic analysis of rocks based on their paleontologic content. Generally speaking, stratigraphy analyses the rock strata and is concerned with their succession and age relationship. All aspects of rocks as strata are, however, of concern for stratigraphy. The analysis of fossil can also provide useful information regarding the environment in which rocks have accumulated: for example, a coral is an unambiguous indication of a warm ocean.

B. The notion of a stratigraphic map

One way of determining the age of a fossil is based on the fact that in a normal sequence the age increases with the depth in the well that penetrates that sequence. So, if the rock accumulation rate is known, the depth x at which the fossil species was found can be used to determine its age y.

The exact dependence between the depth x and the age y – called a *stratigraphic map* – is different for different locations, because it depends on the geological history of this location. The rates of rock accumulation depend of many factors. They vary both in time and space. Thus, in complete geological sequences, fossil taxa of different geologic and absolute ages can be encountered at similar depths in different sections when the section with the younger taxon was accumulated at a higher rate than the one containing the older taxon.

C. Main ideas behind constructing a stratigraphic map

How is the stratigraphic map constructed now? In every area, we have several fossils whose age y has been determined.

The last and the first stratigraphic occurrences of fossil markers occurring in geologic sequences are used for determining the ages of the analyzed samples. Such paleobiological events defined by the same fossil group (e.g., Foraminifera) and indicating different ages may occur in one and the same sample.

- First stratigraphic occurrences are more reliable because such palebiologic events seem to better correlate over extensive areas. They are commonly used when dealing with samples retrieved from outcrops.
- In the analysis of well samples, however, last stratigraphic occurrences are most commonly used because of potential contamination of the deeper samples with material from the overlying rocks.

In both cases, for the selected fossil, we know the depth x_i at which it was found, and we know the estimated age y_i . Based on the points (x_i, y_i) , we must find the desired dependence y = f(x).

Since deeper layers are older, we should have a monotonic (increasing) dependence y = f(x) for which $y_i = f(x_i)$. So, ideally, we should have a monotonic function that passes through all the points.

D. The practical construction of a stratigraphic map is not that easy

The conclusion about monotonicity is based on the idealized assumption that y_i is the age of the oldest (for wells, youngest) of many fossils of this type. For some types, we do have many fossils, so the oldest of these fossils represents a reasonable size sample and is, therefore, highly reliable. For other types of fossils, however, we may have only a few sample fossils of this type in a given area; for such types, the corresponding age y_i and depth x_i are not very accurate.

As a result of this inaccuracy, in practice, it is usually impossible to have a monotonic dependence that passes exactly through all the points (x_i, y_i) : we may have $x_i < x_j$ while $y_i > y_j$.

E. Traditional approach

The traditional paleontological approach to this problem is, crudely speaking, as follows (see, e.g., [8]). Since fewsample data points do not fit to a monotonic curve, we select a threshold n_0 and only consider points (x_i, y_i) which came from samples of size $\geq n_0$.

Ideally, we should select the smallest possible n_0 for which the values can still fit into a monotonic curve, i.e., for which $x_i < x_j$ always implies $y_i \le y_j$.

F. Room for improvement

In the traditional approach, after setting up a threshold, we:

- ignore all the points (x_i, y_i) with lower accuracy, and
- consider all the points with higher accuracy as exact, ignoring the fact that these points are not absolutely accurate.

In both cases, it is desirable to use the ignored information:

- if we take into consideration the inaccuracy of the data (x_i, y_i) based on which we have built the stratigraphic map, then we would be able to determine the accuracy of this map;
- if, in addition to the data points that fit into a monotonic curve, we take into consideration less accurate data points as well, we will be hopefully able to construct a more accurate stratigraphic map.

G. Interval uncertainty

How can we describe the data accuracy? Inaccuracy means that, e.g., for the age, the actual (unknown) age y_i of all the (yet uncovered) fossils of the given type is, in general, somewhat different from the estimate oldest age \tilde{y}_i .

Similarly, because of the additional chaotic rock movements, the ideal depth x_i at which a fossil should be if there were no such movements may be somewhat different from the depth \tilde{x}_i at which the fossil was actually found.

Ideally, we should know the set of possible values of the estimation error $\Delta y_i \stackrel{\text{def}}{=} \widetilde{y}_i - y_i$, and we should know the probabilities of different values from this set. However, to be able to determine these probabilities, we need to have a large number of data points, and when we have a lot of data points, the estimate is pretty accurate anyway. Therefore, in the important cases in which we want to know the accuracy, we cannot experimentally determine these probabilities.

At best, we can find a confidence interval based on the known properties of the extreme-value statistics (see, e.g., [9], [30]), or just elicit these intervals from the experts.

So, if we take uncertainty into consideration, then, for each fossil type *i*, instead of the exact values x_i and y_i , we know the intervals $\mathbf{x}_i = [\underline{x}_i, \overline{x}_i]$ and $\mathbf{y}_i = [\underline{y}_i, \overline{y}_i]$ that contain these unknown values.

H. Fuzzy uncertainty

Often, in addition (or instead) the guaranteed bound for Δy_i , an expert can provide bounds that contain Δy_i with a certain degree of confidence. Usually, we know several such bounding intervals corresponding to different degrees of confidence. Such a nested family of intervals is also called a *fuzzy set*, because it turns out to be equivalent to a more traditional definition of fuzzy set [4], [13], [17], [18], [19] (if a traditional fuzzy set is given, then different intervals from the nested family can be viewed as α -cuts corresponding to different levels of uncertainty α).

I. Towards the precise formulation of the problem

In the case of interval uncertainty, based on all the fossils found in a given area, we know the *n* boxes $\mathbf{x}_i \times \mathbf{y}_i$ corresponding to different types of fossils. We know that the monotonic dependence y = f(x) is such that $y_i = f(x_i)$ for some $(x_i, y_i) \in \mathbf{x}_i \times \mathbf{y}_i$.

Our objective is to find, for every depth x, the bounds of the possible values of age y = f(x) for all the dependencies that are consistent with the given data.

In the case of fuzzy uncertainty, for each degree of confidence α , we must solve the problem corresponding to the α -cut intervals; thus, for each depth x, instead of an interval of possible values of the age y = f(x), we get a nested family of intervals corresponding to different α – i.e., a fuzzy value for the age y = f(x).

These are the problem that we will solve in this paper.

Since technically, the fuzzy problem can be reduced to several interval ones, we will be concentrating on the algorithms for solving the interval problem.

J. Other practical applications of the resulting mathematical problem

Before we find the bounds on f(x), we must first check that our interval bounds are consistent, i.e., that there exists a monotonic function that is consistent with all the boxes. This subproblem has many applications outside paleontology.

Indeed, in many problems in science and engineering, we know that a physical quantity y depends on the physical quantity x, i.e., y = f(x) for some function f(x), and we want to check whether this dependence is monotonic.

In *spectral analysis*, chemical species are identified by locating local maxima of the spectra; see, e.g., [24], [25]. Thus, to identify the chemical species, we must identify intervals between local extrema, i.e., intervals of monotonicity.

In radioastronomy, sources of celestial radio emission and their subcomponents are identified by locating local maxima of the measured brightness of the radio sky. In other words, we are interested in the local maxima of the *brightness* distribution, i.e., of the function y(x) that describes how the intensity y of the signal depends on the position x of the point from which we receive this signal. Thus, in radioastronomy, we must also identify the intervals of monotonicity.

Elementary particles are identified by locating local maxima of the experimental curves that describe (crudely speaking)

the scattering intensity y as a function of energy x. Thus, in elementary particle physics, finding intervals of monotonicity is also important.

In 1-D landscape analysis, e.g., different mountain slopes are different monotonicity intervals; see, e.g., [1], [2], [5].

In *financial analysis*, it is also important to find intervals of monotonicity because they correspond to growth or decline periods; see, e.g., [7]

In *clustering*, different 1-D clusters correspond to a multimodal distribution, so clusters can be naturally described as combinations of monotonicity intervals separating local minima of the probability density function; see, e.g., [15], [21], [22].

Local maxima and minima are also used in the methods that accelerate the convergence of the measurement result to the real value of a physical variable, and thus allow the user to estimate this value without waiting for the oscillations to stop [20]. Thus, to accelerate convergence, we must also be able to efficiently find intervals of monotonicity.

Although checking monotonicity is only a subproblem of the larger problem of locating local extrema, once we know how to efficiently solve this subproblem, we can also efficiently solve the larger problem as well. Indeed, we can find the first interval of monotonicity [1, k] by bisection:

- we originally know that $k \in [k^-, k^+]$, where $k^- = 1$ and $k^+ = n$;
- once we know an interval that contains k, we take its midpoint k_m and check whether $[1, k_m]$ is the interval of monotonicity.
 - If it is, then k ≥ k_m, so we can replace k₋ with k_m.
 If it is not, then k ≤ k_m, so we can replace k⁺ with k_m.

In both cases, we have a half-size interval that is still guaranteed to contain k.

In *l* iterations, we reduce the original width *n* of this interval to $2^{-l} \cdot n$, so in $\log(n)$ iterations, we find *k*. After that, we find the second interval of monotonicity, etc.

K. Additional complexity

Algorithms for solving the subproblem of checking motonoticity have been previously described in [16], [26], [27], [28], [29].

The additional complexity that we face in the biostratigraphy problem comes from the fact that, as we have mentioned, it is possible to have several different ages for the same depth. In mathematical terms, this means that the dependence y = f(x) is not necessarily a monotonic function, it may be a *limit* of the graphs of monotonic functions in the sense of Hausdorff metric (see, e.g., [23]).

We are now ready for the exact definitions and for the formulation of the result.

II. PRECISE FORMULATION OF THE PROBLEM AND THE MAIN RESULT

Definition 1: By a monotonic dependence f, we mean the graph of a continuous mapping $m(s) = (m_1(s), m_2(s))$ from

the real line \mathbb{R} to the plane \mathbb{R}^2 for which t < s implies that $m_1(t) \leq m_1(s)$ and $m_2(t) \leq m_2(s)$.

It is easy to see that if the graph f is the graph of a function, then this definition is equivalent to this function being (non-strictly) monotonically increasing. Not every monotonic dependence is a function: e.g., a "step-function" for which y = 0 for x < 0, y = 1 for x = 1, and $y \in [0, 1]$ for x = 0, is a monotonic dependence but not a function.

Definition 2: By a *box*, we mean a Cartesian product of two intervals. We say that a monotonic dependence f is *consistent* with a box $\mathbf{x} \times \mathbf{y}$ if the graph f contains a point from this box, i.e., if $f \cap (\mathbf{x} \times \mathbf{y}) \neq \emptyset$.

Definition 3: By data d, we mean a finite collection of boxes. We say that a monotonic dependence f is consistent with the data d – and denote it Con(f, d) – if f is consistent with each of the corresponding boxes. We say that the data d is consistent if there exists a monotonic dependence f that is consistent with this data.

Theorem 1: The data d is consistent if and only if for every i and j, $\overline{x}_i < \underline{x}_j$ implies $\underline{y}_i \leq \overline{y}_j$.

For consistent data, our objective is, given the data $[\underline{x}_i, \overline{x}_i] \times [\underline{y}_i, \overline{y}_i]$ $(1 \le i \le n)$ and a real number x, to find the exact lower and upper bounds of the corresponding y over all the monotonic dependences that are consistent with this data:

$$\underline{f}(x) \stackrel{\text{def}}{=} \inf\{y \,|\, \exists f\left((x, y) \in f \,\& \, Con(f, d)\right)\}. \tag{1}$$

$$\overline{f}(x) \stackrel{\text{def}}{=} \sup\{y \,|\, \exists f\left((x, y) \in f \& Con(f, d)\right)\}.$$
(2)

Theorem 2:

$$\underline{f}(x) = \max_{i:\overline{x}_i < x} \underline{y}_i; \quad \overline{f}(x) = \min_{j:x < \underline{x}_j} \overline{y}_j. \tag{3}$$

Proof: Let us denote

$$\underline{F}(x) \stackrel{\text{def}}{=} \max_{i:\overline{x}_i < x} \underline{y}_i; \quad \overline{F}(x) \stackrel{\text{def}}{=} \min_{j:x < \underline{x}_j} \overline{y}_j. \tag{4}$$

Let us first show that for every monotonic dependence f that is consistent with the given data d and for every y for which $(x, y) \in f$, the value y is located between $\underline{F}(x)$ and $\overline{F}(x)$.

Indeed, for every *i*, since *f* is consistent with *d*, there exists a pair $(x_i, y_i) \in f \cap (\mathbf{x}_i \times \mathbf{y}_i)$. For this pair, $\underline{x}_i \leq x_i \leq \overline{x}_i$, so $\overline{x}_i < x$ implies that $x_i < x$. Since $(x, y) \in f$ and $(x_i, y_i) \in f$, by definition of a monotonic dependence, the inequality $x_i < x$ implies that $y_i \leq y$. Since $\underline{y}_i \leq y_i$, we thus conclude that $y \geq y_i$.

Since y is greater than or equal to \underline{y}_i for all *i* for which $\overline{x}_i < x$, it is therefore greater than or equal to the largest of such \underline{y}_i , i.e., that $y \ge \underline{F}(x)$. We can similarly prove that $y \le \overline{F}(x)$.

If $\overline{x}_i < \underline{x}_j$, then, for any x from the open interval $(\overline{x}_i, \underline{x}_j)$, we have $\overline{x}_i < x < \underline{x}_j$. We have proven that for every monotonic dependence that is consistent with the data d, we

have $\underline{y}_i \leq y \leq \overline{y}_j$. So, if the data *d* is consistent, then $\overline{x}_i < \underline{x}_j$ indeed implies $\underline{y}_i \leq \overline{y}_j$ – this is exactly the condition from Theorem 1.

Vice versa, if this condition is satisfied, then we always have $\underline{F}(x) \leq \overline{F}(x)$.

To complete the proof of Theorems 1 and 2, it is therefore sufficient to prove that both piecewise constant monotonic functions $\underline{F}(x)$ and $\overline{F}(x)$, when extended to stepwise monotonic continuous dependences, are consistent with the data d, i.e., that for every k, each of these functions is consistent with the k-th box $\mathbf{x}_k \times \mathbf{y}_k$. Without losing generality, let us prove it for F(x).

Indeed, for a piecewise-constant step dependence like $\underline{F}(x)$, at each point x, the range of possible values of y goes from $\underline{F}(x-0) \stackrel{\text{def}}{=} \lim_{\varepsilon > 0, \varepsilon \to 0} \underline{F}(x-\varepsilon)$ to $\underline{F}(x+0) \stackrel{\text{def}}{=} \lim_{\varepsilon > 0, \varepsilon \to 0} \underline{F}(x+\varepsilon)$. Due to monotonicity, when x goes from \underline{x}_k to \overline{x}_k , possible values of y go from $\underline{F}(\underline{x}_k - 0)$ to $\underline{F}(\overline{x}_k + 0)$. (Since the graph \underline{F} is a graph of a continuous mapping from the real line, it is connected, so all the values from the corresponding intervals are possible.) Therefore, to prove that this graph intersects with the box, it is sufficient to prove that one of the possible values of y also belongs to the y-interval $[y_k, \overline{y}_k]$, i.e., that

$$[\underline{F}(\underline{x}_k - 0), \underline{F}(\overline{x}_k + 0)] \cap [\underline{y}_k, \overline{y}_k] \neq \emptyset.$$
(5)

The formula for the intersection of the two intervals is well known: $[a, b] \cap [a', b'] = [\max(a, a'), \min(b, b')]$. Thus, the two intervals have a non-empty intersection if and only if $\max(a, a') \leq \min(b, b')$, i.e., if and only if $a \leq b'$ and $a' \leq b$. In our case, we must prove that $\underline{F}(\underline{x}_k - 0) \leq \overline{y}_k$ and $\underline{y}_k \leq \underline{F}(\overline{x}_k + 0)$.

By definition of $\underline{F}(x)$ (formula (4), we have:

$$\underline{F}(\overline{x}_k + \varepsilon) = \max_{i:\overline{x}_i < \overline{x}_k + \varepsilon} \underline{y}_i.$$
(6)

Since $\varepsilon > 0$, the inequality $\overline{x}_i < \overline{x}_k + \varepsilon$ holds for i = k. Thus, $\underline{F}(\overline{x}_k + \varepsilon)$ is the largest of several values including \underline{y}_k . Hence, $\underline{F}(\overline{x}_k + \varepsilon) \ge \underline{y}_k$, and in the limit $\varepsilon \to 0$, we get the desired inequality $\underline{F}(\overline{x}_k + 0) \ge \underline{y}_k$.

Similarly,

$$\underline{F}(\underline{x}_k - \varepsilon) = \max_{i:\overline{x}_i < \underline{x}_k - \varepsilon} \underline{y}_i.$$
(7)

The inequality $\overline{x}_i < \underline{x}_k - \varepsilon$ implies that $\overline{x}_i < \underline{x}_k$. We already know that this new inequality, in its turn, implies that $\underline{y}_i \leq \overline{y}_k$. Since all the maximized values \underline{y}_i do not exceed \overline{y}_k , the largest of these values, i.e., $\underline{F}(\underline{x}_k - \varepsilon)$, also cannot exceed \overline{y}_k . In the limit, we get $\underline{F}(\underline{x}_k - 0) \leq \overline{y}_k$.

Both inequalities have been proven, and so are Theorems 1 and 2.

III. RESULTING EFFECTIVE ALGORITHMS

A. Algorithms for checking consistency

For checking consistency, Theorem 1 leads to exactly the same condition as emerged, for a slightly different problem, in [26], [27]. We can therefore use algorithms from [26], [27] to check consistency of our data as well.

If we simply check the condition from Theorem 1 for all i = 1, ..., n and all j = 1, ..., n, then checking this condition would require $O(n^2)$ comparisons – i.e., $O(n^2)$ computational steps.

We can check this condition faster if we use the fact that this condition is equivalent to the following auxiliary property:

For every *i*, we have
$$\underline{y}_i \leq \min_{j:\underline{x}_j \geq \overline{x}_i} \overline{y}_j$$
. (8)

To check this condition, we can perform the following fourstage algorithm:

• First, we sort the values \underline{x}_i into an increasing sequence – this requires $O(n \cdot \log(n))$ steps. We correspondingly re-order the values \overline{x}_i , \underline{y}_i , and \overline{y}_i . After this stage, we can assume that the values \underline{x}_i are sorted:

$$\underline{x}_1 \leq \underline{x}_2 \leq \ldots \leq \underline{x}_n$$

- Then, for every *i* from 1 to *n*, we compute the value $M_i \stackrel{\text{def}}{=} \min(\overline{y}_n, \overline{y}_{n-1}, \dots, \overline{y}_i)$. Here, $M_n = \overline{y}_n$. If we already know M_i , then we can compute the previous value M_{i-1} by using a single operation $M_{i-1} = \min(M_i, \overline{y}_{i-1})$. Thus, computing all *n* values requires *n* computational steps.
- For each *i* from 1 to *n*, we can now use binary search (see, e.g., [6]) to find the integer *m(i)* for which <u>x</u>_{m(i)-1} < *x̄*_i ≤ <u>x</u>_{m(i)} (if such a value exists). Each binary search requires log(*n*) computational steps; thus, *n* such searches require O(n · log(n)) steps.
- Finally, for every i from 1 to n for which m(i) exists, we check whether y_i ≤ M_{m(i)}:
 - if this inequality holds for all such i, then the measurement data is consistent with monotonicity;

- otherwise, the function f(x) cannot be monotonic.

Each checking requires one comparison, so to check that this inequality holds for all i from 1 to n, we need n comparisons.

Overall, we thus need $O(n \cdot \log(n)) + O(n \cdot \log(n)) + O(n) + O(n) = O(n \cdot \log(n))$ steps.

For large n, we may want to further speed up computations if we have several processors working in parallel. This may not be that important for biostratigraphy, but it may be important for other applications of this algorithm. All fours stages of the above algorithm can be parallelized by known techniques. In particular, Stage 2 is a particular case of a general *prefix-sum* problem, in which we must compute the values a_n , $a_n * a_{n-1}$, $a_n * a_{n-1} * a_{n-2}$, ..., for some associative operation * (in our case, * = min).

If we have a potentially unlimited number of processors, then we can do the following (see, e.g., [11], for the information on how to parallelize the corresponding stages):

- on Stage 1, we can sort the values \underline{x}_i in time $O(\log(n))$;
- on Stage 2, we can compute the values M_i (i.e., solve the prefix-sum problem) in time $O(\log(n))$;
- on Stage 3, we can use n processors, each of which compute the corresponding value m(i) in time $O(\log(n))$;

• finally, on Stage 4, we can use n processors, each of which checks the corresponding inequality in time O(1).

As a result, we can check monotonicity in time

$$O(\log(n)) + O(\log(n)) + O(\log(n)) + O(1) = O(\log(n)).$$

If we have p < n processors, then we can:

- on Stage 1, sort n values in time $O((n \cdot \log(n))/p + \log(n))$ [11];
- on Stage 2, compute the values M_i in time $O(n/p + \log(p))$ [3];
- on Stage 3, we subdivide n indices i between p processors, so each processor computes m(i) for n/p indices i; computing each index requires log(n) time, so the overall time is (n/p) · log(n) = O((n · log(n))/p);
- finally, on Stage 4, each of p processors checks the desired inequality for its n/p indices; this requires time O(n/p).

Overall, we thus need time $O\left(\frac{n \cdot \log(n)}{p} + \log(p)\right)$.

B. Algorithms for constructing lower and upper bounds

The function $\underline{f}(x)$ as described by the formula (3) is piecewise constant; when x increases, the value of $\underline{f}(x)$ can only change if when $x = \overline{x}_i$ for some i.

Thus, to compute the corresponding values of $\underline{f}(x)$, it is sufficient to sort the upper endpoints \overline{x}_i into the increasing sequence $\overline{x}_1 \leq \overline{x}_2 \leq \ldots \leq \overline{x}_n$, and then to compute the corresponding values $m_i \stackrel{\text{def}}{=} \max(\underline{y}_1, \ldots, \underline{y}_i)$.

Similarly to the previous algorithm, sorting requires $O(n \cdot \log(n))$ steps and computing m_i requires n steps, so overall, we need $O(n \cdot \log(n))$ steps to compute $\underline{f}(x)$. Similarly, we need $O(n \cdot \log(n))$ steps to compute $\overline{f}(x)$, so the overall computational complexity is $O(n \cdot \log(n))$.

If we have a potentially unlimited number of processors working in parallel, then sorting requires time $O(\log(n))$ and computing m_i also requires time $O(\log(n))$, so overall, we need time $O(\log(n))$. If we have p < n processors, then we need time $O\left(\frac{n \cdot \log(n)}{p} + \log(p)\right)$.

Overall, the computation time for computing the bounds is asymptotically the same as the time for checking consistency.

IV. FROM MONOTONICITY TO MORE COMPLEX CONSTRAINTS

In some practical problems, we know not only that the unknown dependence is monotonic, we also know that its rate of increase cannot be smaller than a certain value c > 0. For example, in paleontology, we may know that the accumulation rate cannot exceed a certain value a; then $\frac{dy}{dx} \ge c \stackrel{\text{def}}{=} 1/a$.

In such situations, we face a slightly different problem: given the data d, check whether there is a dependence y(x) that is consistent with the data and for which, $\frac{dy}{dx} \ge c$ for all points x.

The condition $\frac{dy}{dx} \ge c$ is equivalent to $\frac{dz}{dx} \ge 0$ for a new auxiliary variable $z \stackrel{\text{def}}{=} y - c \cdot x$. In terms of (x, z), the original boxes becomes parallelograms: for $x_i = \underline{x}_i$, we have an interval $[\underline{z}_i^-, \overline{z}_i^-] \stackrel{\text{def}}{=} [\underline{y}_i - c \cdot \underline{x}_i, \overline{y}_i - c \cdot \underline{x}_i]$; for $x_i = \overline{x}_i$, we have an interval $[\underline{z}_i^+, \overline{z}_i^+] \stackrel{\text{def}}{=} [\underline{y}_i - c \cdot \overline{x}_i, \overline{y}_i - c \cdot \overline{x}_i]$. So, we can reformulate the original problem as follows: check whether there exists a monotonic dependence g(x) (= $f(x) - c \cdot x$) that is consistent with all the resulting parallelograms.

Here, for every monotonic dependence g that is consistent with the parallelograms, there exists a point $(x_i, z_i) \in g$ that is inside the parallelogram. Thus, for $(x, z) \in g$, $x < \underline{x}_i$ implies $x < x_i$, hence $z \le z_i \le \overline{z}_i^-$. Similarly, $\overline{x}_i < x$ implies $z \ge \underline{z}_i^+$. Thus, $g(x) \le y \le \overline{g}(x)$, where

$$\underline{g}(x) = \max_{i:\overline{x}_i < x} \underline{z}_i^+; \quad \overline{g}(x) = \min_{j:x < \underline{x}_j} \overline{z}_j^-.$$
(9)

Consistency means that $\underline{g}(x) \leq \overline{g}(x)$ for every x, i.e., that $\overline{x}_i < \underline{x}_j$ implies that $\underline{z}_i^+ \leq \overline{z}_i^-$, i.e., substituting the expressions for z in terms of y that x, that $\underline{y}_i - c \cdot \overline{x}_i \leq \overline{y}_j - c \cdot \underline{x}_j$, hence $\overline{y}_j - \underline{y}_i \geq c \cdot (\underline{x}_j - \overline{x}_i)$, which is equivalent to $c \leq (\overline{y}_j - \underline{y}_i)/(\underline{x}_j - \overline{x}_i)$. Similarly to the monotonic case, one can prove that the above expressions $\underline{g}(x)$ and $\overline{g}(x)$ are indeed the exact bounds on possible values of $z = y - c \cdot x$; thus, $g(x) + c \cdot x$ and $\overline{g}(x) + c \cdot x$ are the exact bounds for y.

A natural next question is: what are the possible values of dy/dx? For every data, we can consider all the "differentiable" functions (in the limit-motivated generalized sense, to allow step functions) that are consistent with all the boxes (i.e., whose graphs intersect with all the boxes). For a given interval [a, b], for each of such functions f, we can take a connected interval hull co(f'([a, b])) of the range of the derivative, and consider the intersection F'([a, b]) of these ranges over all such f.

If this intersection F'([a, b]) contains negative values, this means that every function that is consistent with the data is sometimes decreasing, so no monotonically increasing function is consistent with the data.

In general, the above arguments show, in effect, that the range F'([a,b]) is equal to $\{x \mid p \le x \le q\}$, where

$$p \stackrel{\text{def}}{=} \min_{i,j:a \leq \overline{x}_i \leq \underline{x}_j \leq b} \frac{\underline{y}_j - \overline{y}_i}{\underline{x}_j - \overline{x}_i}, \quad q \stackrel{\text{def}}{=} \min_{i,j:a \leq \overline{x}_i \leq \underline{x}_j \leq b} \frac{\overline{y}_j - \underline{y}_i}{\underline{x}_j - \overline{x}_i}.$$

(So, if p > q, the range is empty.) This formula provides a $O(n^2)$ time algorithm for computing the range.

Comment. A similar algorithm was proposed in [15], for the special case when we know the exact values of x_i .

V. FUTURE WORK: ALGORITHMS

In some cases, we know not only the boxes $\mathbf{x}_i \times \mathbf{y}_i$, but also the probabilities of different values (x_i, y_i) from these boxes. For such cases, it is desirable to find not only the bounds on the stratigraphic map f(x), but also the probabilities of different monotonic dependences within these bounds. In particular, it is desirable to come up with the most probable dependence among all dependences that are consistent with the given data.

VI. FUTURE WORK: APPLICATIONS

As of now, the paper is rather mathematical and abstract. The study presented in the paper is motivated by applications (Biostratigraphy and a lot of others), but so far, we only have algorithms and their complexity analysis.

We are currently working on the actual applications of these algorithms to biostratigraphy.

Since, as we have mentioned, the corresponding problem is relevant for many other application areas, we decided to present our algorithms, to encourage researchers from other application areas to apply these algorithms to their problems as well.

In view of the typically fuzzy character of expert knowledge, we believe that the results from the fuzzy case should be of special interest in various applications areas.

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