THE COMPUTATIONAL COMPLEXITY OF KNOT GENUS AND SPANNING AREA

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ABSTRACT. We show that 3-MANIFOLD KNOT GENUS, the problem of deciding whether a polygonal knot in a closed three-dimensional manifold bounds a surface of genus at most g, is **NP**-complete. We also show that the problem of deciding whether a curve in a PL manifold bounds a surface of area less than a given constant C is **NP**-hard.

1. Introduction

In this paper we investigate the computational complexity of some problems in three-dimensional topology and geometry. We show that 3-MANIFOLD KNOT GENUS, the problem of determining a bound on the genus of a knot in a 3-manifold, is \mathbf{NP} -complete. Using similar ideas, we show that the problem MINI-MAL SPANNING AREA, deciding whether a curve in a metrized PL 3-manifold bounds a surface of area less than a given constant C, is \mathbf{NP} -hard.

Determining whether a given knot is trivial or not is historically one of the central questions in topology. The problem of finding an algorithm to determine knot triviality was posed by Dehn [1]. Dehn's investigations into this area led to the formulation of the word and isomorphism problems, which played an important role in the development of the theory of algorithms. The first algorithm for UNKNOTTING was given by Haken [3]. Haken's procedure is based on normal surface theory, a method of representing surfaces originated by Kneser [14]. Analysis of the computational complexity of this algorithm is more recent. Hass, Lagarias and Pippenger showed that Haken's unknotting algorithm runs in time at most c^t , where the knot K is embedded in the 1-skeleton of a triangulated manifold M with t tetrahedra, and c is a constant independent of M or K [5]. It was also shown in [5] that UNKNOTTING is \mathbf{NP} .

The notion of genus was defined by Seifert [23] in 1935 for knots in the 3-sphere, and extends directly to knots in an arbitrary 3-manifold M. Given a knot K, consider the class S(K) of all orientable spanning surfaces for K. These are surfaces embedded in M with a single boundary component that coincides with K. Seifert showed that this class is non-empty for any knot K in the 3-sphere. For

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knots in a general manifold, S(K) is non-empty when K represents a trivial element in the first integer homology group of M. Any oriented surface without boundary can be obtained from a sphere by adding "handles". The number of handles is the genus of the surface: a sphere has genus 0, a torus has genus 1, etc. The genus is also equal to the number of curves along which a surface can be cut while leaving it connected. If a surface has boundary, its genus is defined as the genus of the surface gotten by gluing a disk to each component of the boundary. The genus of a surface can be computed using any subdivision into polygons, from the formula

Euler characteristic =
$$\#$$
vertices - $\#$ edges + $\#$ faces
= $2 - 2 \times \text{genus} - \#$ boundary curves

The genus g(K) of a knot K is the minimum genus of a surface in $\mathcal{S}(K)$, or ∞ if $\mathcal{S}(K) = \emptyset$. The genus measures one aspect of the degree of "knottedness" of a curve.

UNKNOTTING is a special case of the more general problem 3-MANIFOLD KNOT GENUS. This problem asks for a procedure to determine the *knot genus*, the minimal genus of an orientable spanning surface for a knot in a 3-dimensional manifold. A knot is trivial, or unknotted, precisely when its genus is zero. We will show that this more general problem is **NP**-complete. The only previous results obtained on this problem were given in [5], where it was shown to lie in **PSPACE**. No lower bounds on the running time were previously known.

We work with 3-manifolds that are triangulated and orientable, and with orientable embedded surfaces. This is not a significant restriction, since all compact 3-dimensional manifolds admit unique PL structures [15]. A knot in a triangulated 3-manifold M is a connected simple (non self-intersecting) closed curve in the 1-skeleton of M.

We formulate the problem of computing the genus as a language-recognition problem in the usual way, see [2]. In 1961 Schubert [21], in an extension of Haken's work, showed the decidability of the problem:

Problem: 3-MANIFOLD KNOT GENUS

INSTANCE: A triangulated 3-dimensional manifold M, a knot K in the

1-skeleton of M, and a natural number g.

QUESTION: Does the knot K have q(K) < q?

The size of an instance is given by the sum of the number of tetrahedra in M and $\log g$. In Section 3 we establish

Theorem 1. 3-MANIFOLD KNOT GENUS is NP-hard.

It was established in [5] that 3-MANIFOLD KNOT GENUS is in **PSPACE**. In Section 5 we improve this bound to show that it lies in **NP**.

Theorem 2. 3-MANIFOLD KNOT GENUS is **NP**.

In combination these two results give:

Theorem 3. 3-MANIFOLD KNOT GENUS is **NP**-complete.

Theorem 1 is proved through a construction that transforms an instance of ONE-IN-THREE SAT to an instance of 3-MANIFOLD KNOT GENUS. To a boolean expression we associate a positive integer g and a certain knot in a compact 3-manifold. This knot bounds a surface of genus at most g exactly when there is a

truth assignment to the boolean expression satisfying the requirements of ONE-IN-THREE SAT. Since ONE-IN-THREE SAT is **NP**-hard, this establishes that 3-MANIFOLD KNOT GENUS is also **NP**-hard.

In Section 5 we prove Theorem 2, giving a certificate which demonstrates that a genus g knot K bounds a surface of genus at most g. In [5], which established that UNKNOTTING is \mathbf{NP} , the argument relies on the existence of a normal disk that lies along an extremal ray in the space of normal solutions. The existence of such an extremal normal surface of minimal genus spanning a knot is not known, so a new technique is needed. This is provided by Theorem 12, which gives an algorithm to count the number of orbits of a type of pseudogroup action on a set. Theorem 12 seems likely to have more general applicability. In Section 4 we describe this algorithm and in Section 5 we apply it to a pseudogroup action that arises in the theory of normal surfaces. This allows us to determine in polynomial time the number of components in a normal surface described by an integer vector in \mathbb{Z}_+^{7t} . In particular we are able to certify that a normal surface is connected, orientable and has connected boundary. Since calculating the Euler characteristic of a normal surface is straightforward, establishing orientability and connectedness are the key steps in constructing a polynomial time certificate of its genus.

In Section 6 we extend the orbit counting algorithm to allow the counting of additional integer weight sums associated to each orbit. This allows for the polynomial time calculation of the genus of all the components of a fundamental normal surface, as well as a count of the number of components.

The genus and the area of a surface are closely connected, for example via the Gauss-Bonnet Theorem. In Section 7 we extend the methods developed in studying genus to study the problem of determining the smallest area of a spanning surface for a curve in a 3-manifold. We show that computing an upper bound on the area of a smallest area spanning surface is **NP**-hard.

See [16] for a discussion of complexity classes such as **NP** and **PSPACE** and [25] for a survey of complexity problems in low-dimensional topology.

Remarks:

- 1. Knots are often studied in \mathbb{R}^3 or S^3 rather than in a general manifold. Our methods show that determining knot genus in \mathbb{R}^3 or S^3 , or any fixed manifold, is **NP**. It is not clear whether the problem remains **NP**-hard in \mathbb{R}^3 or S^3 .
- 2. In a related result, Casson has shown that a procedure to determine whether a 3-manifold is homeomorphic to the 3-sphere, following the 3-sphere recognition algorithm described in [18] and [24], runs in time less than $3^t p(t)$, where p(t) is a polynomial. In the direction of lower bounds, it was shown in [13] that determining certain values of the Jones polynomial of alternating links is $\#\mathbf{P}$ -hard.

2. Normal Surfaces

General surfaces in 3-manifolds can wind and twist around the manifold in complicated ways. Kneser described a procedure in which surfaces can be "pulled taut", until they take a simple and rigid position [14]. In this normal position, they have very succinct algebraic descriptions. We use an approach to normal surfaces based on work of Jaco and Rubinstein [12] and Jaco and Tollefson [11]. A normal surface S in a triangulated compact 3-manifold M is a PL-surface whose intersection with

each tetrahedron in M consists of a disjoint set of elementary disks. These are either triangles or quadrilaterals. See Figure 1.

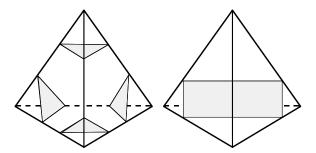


FIGURE 1. Elementary disks in a normal surface.

Within each tetrahedron of M there are four possible triangles and three possible quadrilaterals, up to isotopies that map the tetrahedron to itself. A normal isotopy is an isotopy which leaves each cell of the triangulation invariant. W. Haken observed that a normal surface is determined up to such isotopies by the number of pieces of each of the seven kinds of elementary disks that occur in each tetrahedron, or a vector in \mathbb{Z}_+^{7t} . A normal surface is described by a non-negative integer vector $\mathbf{v} = \mathbf{v}(S) \in \mathbb{Z}^{7t}$, that gives the normal coordinates of S. The set of allowable values $\mathbf{v}(S)$ for a normal surface S lies in a certain homogeneous rational cone \mathcal{C}_M in \mathbb{R}^{7t} , called the Haken normal cone. If $\mathbf{v} = (v_1, v_2, \dots, v_{7t}) \in \mathbb{R}^{7t}$, then the Haken normal cone is specified by linear equations and inequalities of the form

$$v_{i_1}+v_{i_2}=v_{i_3}+v_{i_4}$$
 (up to $6t$ equations) ,
$$v_i\geq 0\ , \quad \text{for } 1\leq i\leq 7t\ .$$

The first set of equations expresses matching conditions, which say that the number of edges on a common triangular face of two adjacent tetrahedra, coming from a collection of elementary disks in each of the tetrahedra, must match. For each triangular face there are three types of edges (specified by a pair of edges on the triangle), which yield 3 matching conditions per face. Triangular faces in the boundary ∂M give no matching equations. The second set are called the *positivity* conditions. The cone \mathcal{C}_M is rational, because the above equations have integer coefficients. We let $\mathcal{C}_M(\mathbb{Z}) = \mathcal{C}_M \cap \mathbb{Z}^{7t}$ denote the set of integral vectors in the cone \mathcal{C}_M . An additional set of conditions, the quadrilateral conditions, is required for an integral vector in the cone to correspond to the normal coordinates of an embedded surface. This condition states that of the three types of quadrilateral found in each tetrahedron, only one can occur in the vector with non-zero coefficient. The quadrilateral conditions are required because two distinct types of quadrilateral in a single tetrahedron necessarily intersect, and we are interested in embedded, non-self-intersecting surfaces. A vector in the Haken normal cone that satisfies the quadrilateral conditions corresponds to an embedded normal surface. This surface is unique up to an isotopy fixing each tetrahedron (called a normal isotopy). However it is important to note that the surface may not be connected.

A fundamental normal surface is a normal surface S such that

$$\mathbf{v}(S) \neq \mathbf{v}_1 + \mathbf{v}_2$$
, with $\mathbf{v}_1, \mathbf{v}_2 \in \mathcal{C}_M(\mathbb{Z}) \setminus \{\mathbf{0}\}$.

In the terminology of integer programming, such a vector $\mathbf{v}(S)$ is an element of the minimal Hilbert basis $\mathbb{H}(\mathcal{C}_M)$ of \mathcal{C}_M , see Schrijver [20, Theorem 16.4]. A fundamental normal surface is always connected, but connected normal surfaces need not be fundamental. Hass, Lagarias and Pippenger [5] gave a bound for the complexity of any fundamental surface.

Theorem 4. Let M be a triangulated compact 3-manifold, possibly with boundary, that contains t tetrahedra.

- Any vertex minimal solution $\mathbf{v} \in \mathbb{Z}^{7t}$ of the Haken normal cone C_M in \mathbb{R}^{7t} has $\max_{1 \leq i \leq 7t} (v_i) \leq 2^{7t-1}$.
- Any minimal Hilbert basis element $\mathbf{v} \in \mathbb{Z}^{7t}$ of the Haken fundamental cone C_M has $\max_{1 \leq i \leq 7t} (v_i) < t \cdot 2^{7t+2}$.

Schubert [21] showed that a surface of smallest genus spanning K can be found among the fundamental surfaces.

Theorem 5. There is a minimal genus spanning surface for K which is a fundamental normal surface.

Similar to the theory of normal surfaces, though somewhat easier, is the theory of normal curves. These are curves on a surface that intersect each triangle in a collection of *normal arcs*, arcs that have endpoints on distinct edges. Since there are three such arcs, up to isotopy, in each triangle, normal curves are described by integer vectors in \mathbb{Z}^{3t}_+ in a surface that contains t triangles. Normal curves arise as the boundaries of normal surfaces in a manifold with boundary.

3. 3-MANIFOLD KNOT GENUS IS **NP**-HARD

In this section we show how to reduce an instance of ONE-IN-THREE SAT to an instance of 3-MANIFOLD KNOT GENUS. Since ONE-IN-THREE SAT is known to be **NP**-hard, this establishes that 3-MANIFOLD KNOT GENUS is also **NP**-hard. The problem ONE-IN-THREE SAT concerns logical expressions involving collections of variables (clauses).

Problem: ONE-IN-THREE SAT

INSTANCE: A set U of variables and a collection C of clauses over U such that each clause $c \in C$ has |c| = 3.

QUESTION: Is there a truth assignment for U such that each clause in C has exactly one true literal?

Schaefer [19] established that ONE-IN-THREE SAT is **NP**-complete. To prove Theorem 1, establishing that 3-MANIFOLD KNOT GENUS is **NP**-hard, we show that an arbitrary problem in ONE-IN-THREE SAT can be reduced in polynomial time to a problem in 3-MANIFOLD KNOT GENUS. See Garey and Johnson [2] for many examples of such reductions.

Let $U = \{u_1, u_2, \dots u_n\}$ be a set of variables and $C = \{c_1, c_2, \dots c_m\}$ be a set of clauses in an arbitrary instance of ONE-IN-THREE SAT. We will describe a knot K in a compact 3-dimensional manifold (with no boundary) and an integer g such that K bounds a surface of genus smaller or equal to g if and only if C is satisfiable so that each clause in C contains exactly one true literal. We construct the 3-manifold M in stages. First we construct a 2-dimensional simplicial complex, then we thicken this complex, replacing triangles with subdivided triangular prisms,

getting a triangulated, 3-dimensional manifold with boundary, as indicated in Figure 2. Finally we use a doubling construction, taking two copies of the manifold with boundary and gluing their boundaries together, to obtain a closed 3-manifold.

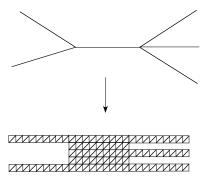


FIGURE 2. A branched surface B, shown in cross-section, is "thick-ened" to produce a triangulated 3-manifold with boundary.

To begin, we form a "branched" surface B by identifying boundary curves of a collection of 2n+1 surfaces with boundary. We construct this collection of surfaces. Let k_i be the number of times that the variable u_i appears in the collection of clauses, and \bar{k}_i be the number of times that \bar{u}_i appears. For $i=1,\ldots,n$, define the surfaces F_{u_i} and $F_{\bar{u}_i}$ to be genus one surfaces with k_i+1 and \bar{k}_i+1 boundary curves respectively. Set F_0 to be a planar surface with n+m+1 boundary curves, one of which will be the knot K. The branched surface B is constructed by identifying these surfaces along appropriate boundary components as indicated in Figure 3.

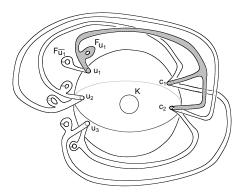


FIGURE 3. A branched surface with boundary curve K corresponding to the boolean expression $(u_1 \vee u_2 \vee u_3) \wedge (u_1 \vee \bar{u}_2 \vee \bar{u}_3)$. The shaded surface F_{u_1} indicates the occurrence of u_1 in each of the clauses c_1 and c_2 .

Branching occurs when more than two boundary curves are identified along a single curve. We identify pairs of boundary curves by giving a homeomorphism between them. Up to isotopy, this is determined by specifying an orientation on the curves and setting the homeomorphism to be orientation reversing. We first fix an orientation on each of F_0 , F_{u_i} and $F_{\bar{u}_i}$. This induces an orientation on each

boundary curve. All identifications will involve gluing a boundary component of F_{u_i} or $F_{\bar{u}_i}$ to a boundary component of F_0 , and we require this gluing to be an orientation reversing homeomorphism. Label by $K, u_1, \ldots u_n, c_1, \ldots c_m$ the 1+n+m boundary components of F_0 . The boundary component K of F_0 , which will become our knot, has nothing identified to it. For each $i, 1 \leq i \leq n$, one boundary curve from the surface F_{u_i} is identified to u_i . The remaining k_i boundary components of F_{u_i} are identified with k_i of the curves $c_1, \ldots c_m$ on ∂F_0 , with one component of ∂F_{u_i} identified with c_j for each occurrence of the literal u_i in the j^{th} clause in C. Similarly, one curve of $\partial F_{\bar{u}_i}$ is identified to u_i , and the remaining \bar{k}_i boundary components are glued to $c_1, \ldots c_m$, with a component glued to c_j for each occurrence of the literal \bar{u}_i in the j^{th} clause of C. A total of three surface boundaries are identified along each of u_1, \ldots, u_n , and exactly four surface boundaries are identified along each of c_1, \ldots, c_m , as in Figure 3.

Lemma 6. There is a truth assignment for U such that each clause in C has exactly one true literal if and only if there is a surface of genus at most m + n mapped continuously into B with boundary K.

Proof. Suppose there is a truth assignment for U such that each clause in C contains exactly one true literal. Form a surface S inside B by taking the union of F_0 and either F_{u_i} if u_i is true, or $F_{\bar{u}_i}$ if u_i is false. Then exactly two boundary components will be identified along each of the boundary components of F_0 other than K itself, and K becomes the boundary of the resulting embedded surface S. There is a contribution of one to the genus from each of the literals, since F_{u_i} and $F_{\bar{u}_i}$ each have genus one, and a contribution of one to the genus from each handle formed when a boundary component of a surface F_{u_i} or $F_{\bar{u}_i}$ is glued to F_0 along a curve c_i . The genus of S is therefore equal to m+n.

Now suppose there is a surface S_1 of genus $\leq n+m$ mapped continuously into B that has a single boundary component mapped homeomorphically to the boundary curve K. We will show in this case that there is a surface S_4 with the same boundary, consisting of pieces of B identified along their boundaries, having genus precisely n+m, and containing, along each curve c_j , exactly one of the three pieces of surface joining F_0 .

The image of S_1 may be quite complicated, winding back and forth across B, but by standard transversality arguments we can homotop S_1 so that it is formed from a union of surface pieces, each piece being mapped homeomorphically to one of the subsurfaces F_0 , F_{u_i} , $F_{\bar{u}_i}$ forming B. More precisely, we can perturb S_1 by a small homotopy so that its intersection with the boundary components of F_0 is transverse, and pulls back to a collection of simple closed curves on S_1 . If any of these curves bounds a disk in S_1 , then the disk is mapped into some subsurface X of B while the boundary curve is mapped to ∂X . Since X is not itself a disk, the disk can be homotoped into ∂X , and we can therefore homotop S_1 in a neighborhood of this disk to remove a component of $S_1 \cap \partial F_0$. After repeating finitely many times, each component of the complement of $S_1 \cap \partial F_0$ in S_1 has non-positive Euler characteristic. The image of S_1 in each subsurface F_0 , F_{u_i} , $F_{\bar{u}_i}$ has an algebraic degree, which is either even or odd. This degree equals the number of pre-images in S_1 of a generic point in the subsurface. The degree of S_1 on F_0 is odd, since K is the boundary of S_1 and therefore S_1 maps an odd number of times to points near K. The sum of the degrees along each of the subsurfaces meeting a curve u_i or c_i is even, since S_1 has no boundary along these curves. In particular, for each $1 \leq i \leq n$, exactly one of F_{u_i} , $F_{\bar{u}_i}$ has odd multiplicity in S_1 . Form a new surface S_2 by taking the union of F_0 and each of the surfaces F_{u_i} , $F_{\bar{u}_i}$ which have odd multiplicity in S_1 . The surface S_2 is obtained from S_1 by a series of operations that either discard a surface with non-positive Euler characteristic or replace a component of S_1 that maps with odd degree to F_{u_i} or $F_{\bar{u}_i}$ with a copy of F_{u_i} or $F_{\bar{u}_i}$. The Euler characteristic of a discarded piece of S_1 is smaller or equal to that of the surface that replaces it, so $\chi(S_2) \geq \chi(S_1)$ in either case.

The collection of subsurfaces of $\{F_{u_i}, F_{\bar{u}_i}\}$ contained in S_2 can be attached to F_0 along common boundary curves among $\{u_1, \dots u_n\}$, forming a connected surface S_3 of genus n. Finally, a connected surface S_4 with boundary K is obtained by identifying pairs of curves in ∂S_3 that are mapped to the boundary components $c_1, \ldots c_m$. The Euler characteristic of S_4 is the same as that of S_3 and S_2 , and therefore greater or equal to that of S_1 . So the genus of S_4 is at most that of S_1 , genus $(S_4) \leq n + m$. Each identification of a pair of surface boundaries along $c_1, \ldots c_m$ contributes one to the genus of S_4 . There is at least one such identification along each of the m curves $\{c_i\}$, though there may be two if each of the three surfaces meeting F_0 along c_j has multiplicity one in S_2 . So identification of curves along $c_1, \ldots c_m$ adds at least m to the genus of S_4 , and it follows that genus $(S_4) \geq$ n+m. Since we have seen that genus(S_4) $\leq n+m$, equality must hold. Equality holds when exactly one of the three surface pieces meeting F_0 along c_i has odd multiplicity for each $1 \leq j \leq m$. We then assign the value "TRUE" to a literal u_i if F_{u_i} is used in S_4 , and the value "FALSE" to u_i if $F_{\bar{u}_i}$ is used in S_4 . This gives a truth assignment to U in which each clause in C has exactly one true literal.

To show that 3-MANIFOLD KNOT GENUS is **NP**-hard we reduce in polynomial time an instance of the **NP**-hard problem ONE-IN-THREE SAT to an instance of 3-MANIFOLD KNOT GENUS.

Proof of Theorem 1. Given an instance of ONE-IN-THREE SAT form B as in Lemma 6. Then there is a truth assignment for U such that each clause in C has exactly one true literal if and only if K is the boundary of a surface of genus m+n mapped continuously into B. Form a 3-manifold N by thickening B so that it is embedded inside a triangulated 3-dimensional manifold with boundary. The thickening process replaces each subsurface forming B by a product of a surface with an interval, and then glues these surfaces together along portions of their boundaries, as indicated in Figure 2. The curve K remains on the boundary of N, and there is a projection map $p:N\to B$ which fixes K. Form a closed manifold M by doubling N along its boundary, namely by taking two copies of N and identifying them along their boundaries by the identity map. Then M admits an involution τ that fixes K, and has quotient N.

To find a triangulation of N, we first describe an explicit triangulation of the branched surface B. An orientable surface with boundary has a triangulation with one vertex on each boundary component and no vertices in the interior. The number of triangles is 4g + 5c - 4, where g is the genus and c is the number of boundary components. We choose such triangulations for each of the 2n + 1 subsurfaces in the branched surface, and we match them together along boundary components to get a triangulation of B. So B has a triangulation in which the number of triangles is linearly bounded in n + m. For the thickened surfaces F_{u_i} and $F_{\overline{u_i}}$, we first form a cell structure of the surface×I by dividing the product into triangles×I. For the thickening of F_0 , we start by doing the same and then go on to divide each interval into five subintervals. Then we glue these thickened surfaces together to get a cell

structure on N. Each cell is this structure is a prism, or triangle×I. We form a closed 3-manifold M by doubling N along its boundary, gluing two copies of N together along their boundaries to obtain a 3-manifold with no boundary.

Finally, we stellar subdivide the cell structure to get a triangulation. Each prism is divided into 14 tetrahedra, by dividing each rectangular face into four triangles by coning to a vertex in the center of each such face, and then coning the 14 triangles of the boundary of the prism to a vertex added to its center. The number of resulting simplices in M is linearly bounded by n+m.

We now check that if K bounds a surface of genus $g \leq m+n$ with interior in $M \setminus K$ then it bounds a surface of the same genus in B. Suppose that F is an embedded surface in M with boundary K. If F does not already lie in N, then perturb it slightly so that the interior of F meets ∂N transversely in a finite number of simple closed curves and arcs. Using the involution τ , reflect the portion of F not in N into N, forming an immersed surface F' lying in N and with the same boundary as F. The interior of F' remains disjoint from K since K is fixed by the involution. The projection p(F') is a surface of genus g mapped into g with boundary g. So if g bounds an embedded surface of genus at most g mapped into g that g in g has a surface of genus at most g mapped into g. The converse was shown in the proof of Lemma 3. it then follows from Lemma 3 that g bounds an embedded surface of genus at most g mapped into g is satisfiable with each clause containing exactly one true literal.

The construction of B, N and M described above each requires only a linear number of steps in the size of the instance of ONE-IN-THREE SAT with which we started, so that the reduction requires polynomial time.

4. Orbits of interval isometries

In this section we develop a combinatorial procedure that computes the number of orbits of a collection of k isometries between subintervals of an interval $[1,N]\subset\mathbb{Z}$ in time polynomial in $k\log N$. By lining up the intersections of a normal surface with the edges of a triangulation, we obtain such subintervals. Arcs of the normal surface on the faces of the triangulation give rise to correspondences of these intersection points which are subinterval isometries. We can then apply the algorithm developed here to count the number of components of a normal surface.

Assume that we have a set of integers $\{1,2,\ldots,N\}$ and a collection of monotone bijections, $g_i:[a_i,b_i]\to[c_i,d_i],\ 1\leq i\leq k,\ a_i\leq c_i,$ either increasing or decreasing, that are called pairings. If a pairing identifies two intervals [a,b] and [c,d] by sending a to c and b to d, we call it an orientation preserving pairing, and if it sends a to d and d to d, we call it orientation reversing. If d and d if it sends d if it sends d if d

The interval $[1, \ldots, N]$ is divided into equivalence classes by the action of the pairings, which are called *orbits*. We are interested in the orbit structure of the collection, since with appropriate interpretation an orbit corresponds to a connected component of a normal surface. We introduce several simplification processes on the set of pairings in order to analyze the structure of the set of orbits. The following lemma describes the orbits of an orientation preserving pairing whose domain and range together form a connected interval.

Lemma 7. When an orientation preserving pairing $g : [a,b] \to [c,d]$ has $a < c \le b+1$, then the pairing has t = c-a orbits on [a,d].

Proof. Each point in [a, b] greater or equal to c lies in the range of g and can be mapped to a smaller point in [a, b] by a power of g^{-1} . So each orbit on [a, d] has a representative in [a, c-1]. Since the congruence class modulo t of a point is preserved by g, each of the t integers in [a, c-1] lies in a distinct orbit. These points uniquely represent the t orbits.

We say that such a pairing is periodic with period t, and the combined interval [a,d] is a periodic interval of period t. We now show how to merge two pairings with sufficient overlap into a single pairing with the same orbits.

Lemma 8. Let R_1 be a periodic interval with pairing g_1 of period t_1 , and suppose that g_2 is an orientation preserving pairing with translation length t_2 which pairs two distinct intervals of width t_1 in R_1 . Then the orbits of $g_1 \cup g_2$ on R_1 are the same as those of a single periodic action on R_1 of period $GCD(t_1, t_2)$.

Proof. Let J_1 be an interval in R_1 of width t_1 that is in the domain of g_2 and which is paired by g_2 to an interval in R_1 . Each point in R_1 has a unique orbit representative in J_1 under the action of g_1 on R_1 . The interval $g_1(J_1)$ lies in R_1 by assumption. For $x \in J_1$ let f(x) be the unique point in J_1 obtained by carrying $g_2(x)$ back to a point in J_1 by a power of g_1 . The effect of f(x) on J_1 is a shift of t_2 (mod t_1). The orbits of f on J_1 divide the points of J_1 into congruence classes modulo $GCD(t_1, t_2)$. Neither g_1 nor g_2 change the congruence class of a point mod $(GCD(t_1, t_2))$, so a subinterval of width $GCD(t_1, t_2)$ in R_1 contains exactly one representative of each orbit of R_1 under the action of $g_1 \cup g_2$. The same orbits arise from a periodic action of period $GCD(t_1, t_2)$.

The following is a special case of Lemma 8 that applies when both g_1 and g_2 are periodic pairings.

Lemma 9. Let R_1 and R_2 be overlapping periodic intervals, associated with pairings g_1, g_2 having periods t_1 and t_2 . Suppose that width $(R_1 \cap R_2) \ge t_1 + t_2$. Then the orbits of $g_1 \cup g_2$ on $R_1 \cup R_2$ are the same as those of a single periodic pairing on $R_1 \cup R_2$ of period $GCD(t_1, t_2)$.

Proof. The leftmost interval J_1 of width t_1 in $R_1 \cup R_2$ is translated by g_2 a distance of t_2 to the right, to an interval J_2 which lies in $R_1 \cup R_2$. The result follows from Lemma 8.

We state a consequence in a form which will be convenient for our applications.

Lemma 10. Let g_1, g_2 be periodic pairings with periods t_1, t_2 and let J_1, J, J_2 be intervals with $g_1(J_1) = J$ and $g_2(J) = J_2$. Suppose that J is contained in the union $J_1 \cup J_2$. Then the orbits of $g_1 \cup g_2$ are the same as those of a single periodic pairing of period $GCD(t_1, t_2)$, acting on the union of the periodic intervals of g_1, g_2 .

Proof. Let $J_1 = [a_1, b_1], J = [a, b]$ and $J_2 = [a_2, b_2]$. We have $a_2 \ge a_1$ and $J \subset J_1 \cup J_2$. Since $J_1 \cup J_2$ is a connected interval, we have that $a_2 \le b_1 + 1$. Note that $t_1 = b - b_1 = a - a_1$ and $t_2 = b_2 - b = a_2 - a$. Then $t_1 + t_2 = (b - b_1) + (a_2 - a) = (b - a + 1) + (a_2 - b_1 - 1)$. The width of J is (b - a + 1) and we have seen above that $(a_2 - b_1 - 1) \le 0$. So the width of J is at least $t_1 + t_2$ and Lemma 9 implies the conclusion of the lemma.

The orbit counting algorithm applies a series of modifications to the pairings, which we now describe. The operation of replacing g_1 and g_2 by a single periodic action on $R_1 \cup R_2$ of period $GCD(t_1, t_2)$, as in Lemma 10, is called a *periodic merger*. An interval is called *static* if it is in neither the domain nor the range of any pairing, so that its points are identified to no other points by pairings. The operation of *contraction* is performed on a static interval [r, s]. We eliminate this interval, replace [1, N] by [1, N - (s - r + 1)], and alter each g_j by replacing any point x in a domain or range which lies entirely to the right of s by x - (s - r + 1). (Note that this operation leads to an increase of s - r + 1 in the count of the number of orbits, since the eliminated points are each unique representatives of an orbit.)

The next operation simplifies an orientation reversing pairing whose domain and range overlap. Suppose that $g:[a,b] \to [c,d]$ is a pairing with g(a)=d, g(b)=c and $b \ge c$. Define a new pairing $g':[a,(a+d)/2) \to ((a+d)/2,d]$ by restricting the domain and range of g, and say that g' is obtained from g by trimming. The domain and range of a trimmed pairing are disjoint.

If there is an interval in the domain and range of exactly one pairing, then the interval can be "peeled off" without changing the orbit structure, in a way we will describe below. We will apply this operation only to strip off points from the right of the interval [1,N]. When there is a pairing $g:[a,b]\to [c,N]$ and a value N' with $c\le N'+1\le N$, such that all points in the interval [N'+1,N] are in the domain or range of no pairing other than g, then we can perform an operation called truncation of g. Truncation shortens the interval [1,N] to the interval [1,N'], and similarly shortens the domain and range of g. In the orientation preserving case, pairings other than g are unchanged, while g is either eliminated entirely, if c=N'+1, or otherwise replaced by a shortened pairing $g':[a,b-(N-N')]\to [c,N']$. The inverse pairing g^{-1} is correspondingly altered. When there is no ambiguity, we continue to call the shortened pairing by the same name. Note that we can perform this operation when the interval [N'+1,N] intersects both the range and the domain of g.

In the orientation reversing case, we truncate only when g has disjoint domain and image, and [N'+1,N] is contained in the image. Suppose $g:[a,b] \to [c,N]$ is a pairing with g(a) = N, g(b) = c and b < c. If [N'+1,N] is disjoint from the domains and ranges of all pairings other then g and $N'+1 \ge c$ then we eliminate g if N'+1=c, and otherwise replace g by a shortened orientation reversing pairing formed by restricting its domain to [a+N-N',b] and its range to [c,N'].

Finally we define an operation in which two pairings are composed, with one pairing g_1 used to shift down the domain and range of a second pairing g_2 as much as possible. We will use this operation to allow us to peel off pairings from the right of the interval [1, N'] by the previous operation of truncation. Define a maximal pairing to be one whose range contains both N' and the range of any other pairing containing N'. More precisely, define a linear order on pairings using the lexicographical order $(d_i, -c_i, -a_i, -\text{orientation})$, so that the maximal pairing has

the highest upper endpoint, and among those the widest range, and among those the biggest translation distance (if orientable).

As a first step, if g_1 is orientation reversing and has overlapping domain and image, trim g_1 . Now consider an arbitrary second pairing g_2 whose range is contained in the range of g_1 . If the domain of g_2 is not contained in the range of g_1 , form the composite map $g_2' = g_1^{-r} \circ g_2$, where r = 1 if g_1 is orientation reversing and otherwise $r \geq 1$ is the largest integer such that $g_1^{-r+1}([c_2, d_2])$ is contained in the range of g_1 . If the domain of g_2 is also contained in the range of g_1 then form the composite map $g_2' = g_1^{-r} \circ g_2 \circ g_1^s : g_1^{-s}([a_2, b_2]) \to g_1^{-r}([c_2, d_2])$, where r is as above, s = 1 if g_1 is orientation reversing, and otherwise $s \geq 1$ is the largest integer such that $g_1^{-s+1}([a_2, b_2])$ is contained in the range of g_1 . The process of replacing g_2 by $g_1^{-r} \circ g_2 \circ g_1^s$ is called a transmission of g_2 by g_1 .

Lemma 11. Altering a collection of pairings by the operations of periodic merger, trimming, truncation and transmission preserves the number of orbits of the collection. A contraction decreases the number of orbits by the width of the contracted interval.

Proof. We refer to the collection of pairings $g_1, g_2, \dots g_k$ as G and to the new collection of pairings formed by some operation as G'.

We first consider the effect of a contraction, which cuts down the interval [1, N] by removing points which are fixed by the collection of pairings. The number of orbits removed equals the width of the contracted interval.

Consider now the effect of a periodic merger that merges two periodic intervals R_1 and R_2 . Lemma 9 shows that the orbits in $R_1 \cup R_2$ of $g_1 \cup g_2$ are the same as those produced by a single periodic pairing g'. Suppose that x and y are in the same orbit of the action of $G = \{g_1, g_2, \ldots g_k\}$. Then $h \cdot x = y$, where h is some finite word in the elements of G and their inverses. Wherever a g_1 or g_2 occurs in this word we can replace it by a power of g', since g' translates by an amount that divides the translation distance of g_1 and g_2 . So x and y are in the same orbit of $G' = \{g', g'_3, \ldots g'_k\}$. Conversely suppose that that x and y are in the same orbit under G'. Then $h' \cdot x = y$, where h' is some finite word in $g', g'_3, g'_4 \ldots g'_k$. Lemma 9 implies that wherever a g' occurs in h' we can replace it by some word in g_1 and g_2 , since the orbits of g' and $g_1 \cup g_2$ coincide on the periodic interval of g'. So periodic mergers preserve orbits.

Suppose next that $g_i: [a_i, b_i] \to [c_i, d_i]$ is an orientation reversing pairing with $g_i(a_i) = d_i, g(b_i) = c_i \leq c_i$, and that $g_i': [a_i, (a_i + d_i)/2) \to ((a_i + d_i)/2), d_i]$ is obtained by trimming g_i . If x and y are in the same orbit of $G = \{g_1, g_2, \ldots g_k\}$, then there is a sequence of points $x = x_1, x_2, \ldots, x_r = y$ where each x_j is the image of x_{j-1} under a pairing in G. We can replace an occurrence of g_i or g_i^{-1} by g' if the point x_{j-1} is smaller than $(a_i + d_i)/2$ and by g'^{-1} otherwise. So x and y remain in the same orbit under the action of G'. Conversely if x and y are in the same orbit under G' then replacing each occurrence of g' by g and of g'^{-1} by g^{-1} gives a word in G taking x to y. We conclude that trimming preserves orbits.

Next consider the effect of a truncation. Suppose that all points in [N'+1, N] are in the range of exactly one pairing $g:[a,b]\to [c,N]$, and we truncate, shortening to an interval [1,N'] and either eliminating g or shortening it to $g':[a,b-(N-N')]\to [c,N']$ in the orientation preserving case, or to $g':[a+N-N',b]\to [c,N']$ in the orientation reversing case. Suppose that $g=h\cdot x$ where g=h is a reduced product of pairings (a product that does not contain a subproduct of the form g=h: We

can assume that y>x, as otherwise we can replace h by its inverse. Let h' be obtained from h by replacing all occurrences of g with g'. The successive images of x under the subwords of h can never enter and leave the interval [N'+1,N], since they can only enter it under the action of some positive power of g, and only leave it under a negative power of g. These don't occur in succession in a reduced product. The image of a point z under g_j is unchanged unless $g_j=g$, and $g(z)\geq N'+1$. So the image of x under h is the same as its image under h' unless y>N'. In that case h is of the form $h=g_i^Th_1$ where $h_1\cdot x\leq N'$, $r\geq 1$, and $y=g_i^Th_1\cdot x>N'$. So two points in [1,N'] are in the same orbit under G' if and only if they are in the same orbit under G. But each point in [N'+1,N] is in the same G orbit as a point in [1,N']. It follows that the number of orbits is unchanged by truncation.

Finally suppose that a transmission of g_j by g_i replaces the pairing g_j in G by $g'_j = g_i^{-r} \circ g_j \circ g_i^s$. If $y = h \cdot x$ where h is a word in G, then $y = h' \cdot x$ where h' is obtained by replacing every occurrence of g_j with $g_i^k \circ g_j \circ g_i^{-s}$. Similarly if $y = h' \cdot x$ where h' is a word in G', then $y = h \cdot x$ where h is obtained by replacing every occurrence of g'_j with $g_i^{-k} \circ g_j \circ g_i^s$. So transmissions preserve the collection of orbits and don't change their number.

We now describe an algorithm which uses these operations to count the number of orbits of a pseudogroup of pairings acting on [1,N]. We initially set a counter for the number of orbits to zero. A pairing g_i is said to be maximal if $d_i = N$ and $[c_i, N]$ contains the range of any other pairing with an endpoint at N. Let $S = \{1, 2, ..., N\}$ and let $g_i : [a_i, b_i] \to [c_i, d_i]$, $1 \le i \le k$ be a collection of pairings between subintervals of S. The algorithm will contract S and reduce the number of bijections. Denote by N' the current size of the interval as we proceed. The algorithm repeats the following steps, reducing N' and k, until there are no points remaining.

Orbit counting algorithm:

- 1. Delete any pairings that are restrictions of the identity.
- 2. Make any possible contractions and increment the orbit count by the sum of the number of points deleted by the contractions.
- 3. Trim all orientation reversing pairings whose domain and range overlap.
- 4. Find a maximal pairing g_i . Recall that its range contains both N' and the range of any other pairing containing N'. If g_i is periodic, search for other periodic pairings g_j that also have range $[c_j, N']$ containing N'. If any such g_j exists, and if $t_i \leq w_j$, then perform a merger as in Lemma 10, replacing g_i and g_j by a single periodic pairing, with translation distance $GCD(t_i, t_j)$, acting on the union of the domains and ranges of g_i and g_j .
- 5. Find a maximal g_i . For each g_j with $j \not\models i$, if the range of g_j is contained in $[c_i, N']$, transmit g_j by g_i .
- 6. Find the smallest value of c such that the interval [c, N'] intersects the domain and range of exactly one pairing. Truncate the pairing whose range contains the interval [c, N'].
- 7. If the width of the interval [1, N'] is greater than zero, start again with Step (1). Otherwise output the number of orbits and stop.

Theorem 12. The orbit counting algorithm gives the number of orbits of the action of the pairings $\{g_i\}_{i=1}^k$ on $[1, \ldots, N]$ in time bounded by a polynomial in $k \log N$.

Proof. We first check that the orbit counting algorithm correctly counts the number of orbits. In Step (1), deleting a pairing which is the identity on its domain does not change the number of orbits. In Step (2), contracting a static interval removes a number of points that are unique orbit representatives, and these are added to the running total of the orbit count. If all the points have been removed, then there are no more orbits to count and the orbit count is complete. Lemma 11 shows that the operations of transmission, trimming, truncation and merger occurring in Steps (3)-(6) do not change the number of orbits of the series of collections of pairings. The intervals in Step (4) satisfy the hypothesis of Lemma 10, and can be merged without changing the orbit structure of the collection of pairings. Specifically, consider the three intervals J, J_1 and J_2 , where J_1 and J_2 are the domain and range of g_j and J is the range of J_2 under g_i^{-1} . Since g_j is periodic we know that $J_1 \cup J_2$ is a connected interval. We are assuming that $t_i \leq w_j$, so $J \cup J_2$ is also a connected interval. If $t_i \leq t_j$ then J lies in the interval $J_1 \cup J_2$. If $t_j \leq t_i$ then J_1 lies in the interval $J \cup J_2$. In either case Lemma 10 shows that g_i and g_j can be merged. A merger preserves the orbit structure while reducing the number of pairings by one, replacing g_i and g_j by a single periodic pairing acting on the union of their periodic intervals.

In each cycle through these steps the interval width decreases by at least one. Therefore the algorithm terminates, yielding a count of the number of orbits after a finite number of steps. However we will obtain a much better bound on the running time. To do so, we define a complexity which decreases as we iterate the above steps. Recall that $w_i = b_i - a_i + 1$, $1 \le i \le k$ and k is the number of pairings. The complexity X is defined to be

$$X = 4^k \prod_{i=1}^k w_i.$$

Call the process of executing Steps (1)-(7) a *cycle*. We will show that when we run through 2k cycles, X is reduced by a factor of at least two. See the remark at the end of Section 4 for a geometric interpretation of this complexity.

Call an interval [x,y] a Z-active interval if it is the domain or range of a pairing of subintervals of [1,Z] and if $y-x+1 \geq (Z-x+1)/2$, or equivalently, if $y \geq Z/2+x/2-1/2$. Being Z-active corresponds to being relatively close to Z; there is no room for another interval of the same size to the right of y. The value of Z is initially set to N, and as the algorithm proceeds it is reset to the current interval width N' each time the number of pairings k is decreased. The number of pairings decreases when a merger occurs, or when a pairing is truncated to zero width, or when a pairing is transmitted to become trivial (a restriction of the identity) and then contracted.

Claim 1. The union of two Z-active intervals of equal width is a connected interval.

Proof. Suppose that [a,b] and [c,d] are equal width Z-active intervals with $c \ge a$. Then b-a=d-c, $b-a+1 \ge (Z-a+1)/2$ and $d-c+1 \ge (Z-c+1)/2$, so $b \ge Z/2+a/2-1/2 \ge d/2+a/2-1/2=b/2+c/2-1/2$ implying that $b \ge c-1$ and $[a,b] \cup [c,d]$ is connected.

Claim 2. Suppose that the domain [a,b] of a pairing g is not Z-active. Then the image [x',y'] of an interval [x,y] under g^{-1} is not Z-active.

Proof. Suppose that [x', y'] is the image of [x, y] under g^{-1} , where $g : [a, b] \to [c, d]$. By assumption we have that b < Z/2 + a/2 - 1/2. The interval [x', y'] is contained in [a, b] so $a \le x' \le y' \le b$. Then $y' \le Z/2 + a/2 - 1/2 \le Z/2 + x'/2 - 1/2$ and [x', y'] is not Z-active.

Claim 3. After a series of three cycles involving the truncation of three successive maximal pairings g_1, g_2, g_3 , either the number of Z-active intervals decreases or the number of pairings decreases.

Proof. Suppose that a series of three successive maximal pairings g_1, g_2, g_3 occurs in the orbit counting algorithm with no decrease in the number of pairings. When g_1 stops being maximal, it is transmitted by the new maximal pairing g_2 , which is then itself truncated until no longer maximal. It in turn is then transmitted by g_3 . When g_{i+1} transmits g_i , the pairing of $[a_i, b_i]$ with $[c_i, d_i]$ is replaced by a pairing of $[a'_i, b'_i]$ with $[c'_i, d'_i]$ and one of the following three cases occurs.

- 1. $[c_i, d_i]$ is transmitted to a Z-inactive interval by g_{i+1} .
- 2. $[c_i, d_i]$ is transmitted to a Z-active interval $[c'_i, d'_i]$ by g_{i+1} , and $[a'_i, b'_i]$ is Z-active.
- 3. $[c_i, d_i]$ is transmitted to a Z-active interval $[c'_i, d'_i]$ by g_{i+1} and $[a'_i, b'_i]$ is not Z-active.

Since g_1 is initially maximal, its range is initially Z-active. Truncation reduces the range to $[c_1, d'_1]$, at which point the pairing g_2 becomes maximal. We can assume $[c_1, d'_1]$ is still Z-active, or we are done. The interval $[c_1, d'_1]$ is then transmitted by g_2 to an interval $[e_1, f_1]$. In Case (1) this new interval is Z-inactive and the number of Z-active intervals has decreased. In Case (2) the three intervals $[c_1, d'_1]$, $[e_1, f_1]$ and $[a_1, b'_1]$ are all Z-active and of equal width, and therefore satisfy the hypothesis of Claim 1. Lemma 10 then implies that two pairings can be merged, and the number of pairings has decreased. If Case (3) occurs we truncate the maximal pairing g_2 . As we do so we can assume its domain $[a_2, b'_2]$ remains Z-active, or we are done. When the next pairing g_3 becomes maximal and in turn transmits g_2 , we fall into one of the first two cases, and then either have reduced the number of Z-active intervals or can merge two pairings.

Claim 4. Suppose that Z is set to the current interval size N' and that after a series of truncations in which the maximal pairings are successively g_1, g_2, \ldots, g_r , no Z-active intervals remain. Then the complexity X is reduced by a factor of at least two.

Proof. Truncation of a maximal pairing g_i results in its range $[c_i, d_i]$, of width w_i , being reduced to a shorter interval $[c_i, d'_i]$ of width w'_i . This reduces the complexity X by a factor of

$$\frac{w_i'}{w_i} = \frac{d_i' - c_i + 1}{d_i - c_i + 1}.$$

The maximal pairing switches after a truncation if the range of g'_i has truncated sufficiently so that it is contained in the range of g_{i+1} , so that $c_{i+1} \leq c'_i$ and $d_{i+1} = d'_i$. Define d''_i to be the integer that satisfies the equation

$$d_i'' = c_i - 1 + (w_{i+1}) \frac{Z - c_i + 1}{Z - c_{i+1} + 1}.$$

and define

$$w_i'' = d_i'' - c_i + 1.$$

Note that the function

$$f(x) = \frac{x - c_i + 1}{x - c_{i+1} + 1}$$

is increasing with x for $x \ge c_i$, since $c_i \ge c_{i+1}$. So

$$w'_{i} = d'_{i} - c_{i} + 1 = d_{i+1} - c_{i} + 1 = (w_{i+1}) \frac{(d_{i+1} - c_{i} + 1)}{w_{i+1}}$$

$$\leq (w_{i+1}) \frac{Z - c_{i+1} + 1}{Z - c_{i+1} + 1} = d''_{i} - c_{i} + 1 = w''_{i}.$$

After a series of truncations of the pairings g_1, g_2, \ldots, g_r , the complexity X is multiplied by a factor of

Now $Z = d_1$, since g_1 was the first maximal pairing truncated, so $Z - c_1 + 1 = d_1 - c_1 + 1 = w_1$. By assumption, there are no Z-active intervals remaining after g_r is truncated, so that $[c_r, d'_r]$ is Z-inactive. We then have by the definition of Z-active that

$$\left(\frac{Z-c_1+1}{Z-c_r+1}\right)\left(\frac{w_r'}{w_1}\right) = \frac{w_r'}{Z-c_r+1} = \frac{d_r'-c_r+1}{Z-c_r+1} < 1/2.$$

It follows that X is multiplied by a factor smaller than 1/2.

Proof of Theorem 12: The operations involved in one cycle consist of comparisons, additions, subtractions and computing greatest common divisors of the k pairings of G. The number of these operations occurring in each cycle is at most linear in $k \log N$. Pairings are described by pairs of intervals, whose boundary points are integers of size at most N. So the running time of each cycle is polynomial in $k \log N$.

If the number of pairings k decreases during a cycle, then a pairing has been eliminated because its width has truncated to zero, because it has been transmitted and become the identity on its domain, or because two pairings have merged. When two pairings g_1, g_2 merge, the value of k decreases by one. The product $(w_1)(w_2)$ which occurs in X is replaced by the width of the new pairing, which is at most $(w_1 + w_2) \leq 2w_1w_2$. Therefore $X = 4^k \prod_{i=1}^k w_i$ decreases by a factor of at least two after a merger.

Since each pairing has width at most N, the initial complexity is bounded above by $(N)^k 4^k$. By Claim 3, each time we run through three cycles either the number of pairings decreases or the number of Z-active intervals decreases. There are at most k Z-active intervals initially, so after 3k cycles either there is a reduction in

the number of pairings or there are no Z-active intervals remaining. In the second case the complexity X has decreased by a factor of at least two, by Claim 4.

So X decreases by a factor of at least two after 3k cycles, and the complexity reduces to zero after at most $k(2 + \log N)$ successive series of 3k cycles, or after at most $3k^2(2 + \log N)$ cycles. Each cycle runs in time polynomial in $k \log N$, so the total running time is also polynomial in $k \log N$.

We apply Theorem 12 to count the number of components of a normal curve or normal surface. An obvious algorithm to count components proceeds by marking vertices connected by common edges until all vertices in a component are reached. This procedure takes time linear in the number of edges of the curve, which is equal to the sum of the normal coordinates (W below), but we can achieve an exponential improvement. We first look at normal curves.

Corollary 13. Let F be a surface with a triangulation T containing t triangles and let γ be a normal curve in F with normal coordinates summing to W. There is a procedure for counting the number of components of γ which runs in time polynomial in $t \log W$.

Proof. The 1-skeleton of T contains e edges, where $e \leq 3t$. Fix once and for all an ordering of these edges. A normal curve γ intersects each edge of T in a finite number of points. Set N to be the weight W of γ , which is the sum of the number of intersection points of γ with all the edges of T. Label the intersections of γ and the first edge of the 1-skeleton by the integers $1, 2, \ldots, i_1$, the intersections of γ and the jth edge by $i_{j-1}+1, i_{j-1}+2, \ldots, i_j$, and the intersection of γ and the e^{th} edge of T by $i_{e-1}+1, i_{e-1}+2, \ldots, i_e$. Then $i_e=N$. Each triangular face of T has three sets of arcs pairing points of [1, N], with one set running between each pair of edges of the face. To each set of arcs we associate a pairing between the intervals at either end of the arcs, as in Figure 4. All the pairings are orientation reversing in this example. In general some will be orientation preserving, as the edge orientations on any triangle can be arbitrary.

The number of connected components of γ is the same as the number of orbits of an action of a collection of pairings on $\gamma \cap T^{(1)}$, where two points are paired if they are connected by an edge of γ lying in a triangle. This is precisely the number returned by the orbit counting algorithm. The number of pairings is at most 3t and the interval size is W. Applying Theorem 12, we can determine the number of components of the normal curve γ in time polynomial in $t \log W$.

A similar argument applies to normal surfaces.

Corollary 14. Let M be a 3-manifold with a triangulation T containing t tetrahedra and let F be a normal surface in M whose normal coordinates sum to W. There is a procedure for counting the number of components of F which runs in time polynomial in $t \log W$.

Proof. The 2-skeleton of T contains at most 4t faces and the 1-skeleton contains $e \leq 6t$ edges. Fix once and for all an ordering of the edges. A normal surface F intersects each edge in a finite number of points. Set N to equal the weight W of F. As before, order the edges of $T^{(1)}$ in an arbitrary way and label the intersections of F and the jth edge of $T^{(1)}$ by $i_{j-1}+1, i_{j-1}+2, \ldots, i_j$. Again we have $i_e=N$. To a pair of edges on a triangular face of the 2-skeleton we associate a pairing between the intervals at either end of the corresponding set of arcs. There are at most three pairings for each face, and the number of faces is at most 4t, so the number of

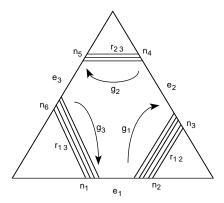


FIGURE 4. Normal arcs on a triangular face give three pairings of ordered intervals: $g_1:[n_1+r_{13},n_2]\to[n_3,n_3+r_{12}-1], g_2:[n_3+r_{12},n_4]\to[n_5,n_5+r_{23}-1],$ and $g_3:[n_5+r_{23},n_6]\to[n_1,n_1+r_{13}-1]$ The normal curve coordinates associated to this face are r_{12},r_{13} and r_{23} . In a normal surface, r_{12},r_{13} and r_{23} are each a sum of two normal coordinates.

pairings is bounded above by 12t. These pairings are determined by the normal coordinates of F.

For a normal surface, the number of connected components of F is the same as the number of components of $F \cap T^{(2)}$, since every component of F intersects the 1-skeleton of T, and if two points on $F \cap T^{(2)} \cap T^{(1)}$ can be joined by a path then that path can be homotoped into the 2-skeleton of the triangulation. So the the number of components equals the number of orbits of $F \cap T^{(1)}$ under pairings that identify two points connected by an edge of $F \cap T^{(2)}$ contained in a face of the 2-skeleton. This number is precisely what is computed by the orbit counting algorithm. Therefore we can determine the number of components of the normal surface F in time polynomial in $t \log W$.

Remark: There is a motivating geometrical construction behind the combinatorics of the orbit counting algorithm which we informally explain. We can associate to each pairing $g_i : [a_i, b_i] \to [c_i, d_i]$ a pair of "transmission towers" in the upper halfplane, with one tower being the vertical line segment from $(a_i, 1/e)$ to (a_i, w_i) and the other the vertical segment from $(c_i, 1/e)$ to (c_i, w_i) . These towers capture the information contained in the pairings. Points in the domain and image of a pairing beam up to the tower leftward at a 45 degree angle, then beam to the paired tower (either straight across, if the pairing preserves orientation, or crossing if not) then down again. Assign a "cost" to each transmission tower equal to its hyperbolic length in the upper half space model. This equals $\log(d_i - c_i + 1) + 1$, and the sum of all these lengths is essentially the logarithm of the complexity X used above.

The counting algorithm starts with a Euclidean line emerging from the right endpoint of an active interval and going upward at a 45 degree angle. In hyperbolic terms, this is an equidistant curve from the geodesic from the right endpoint to the point at infinity. Initially, we may assume that this sweep line hits the top of at least one transmission tower. We make the highest of these the active tower, and use it to beam all the other towers to its paired interval. If the two domains overlap,

we use the highest power of the transmission that can be applied. Eventually the triangle between the current tower and the sweep line is vacant. We sweep leftward with this equidistant curve. As long as the sweep line hits only one tower, we truncate it.

Consider a second equidistant curve P through a point x on \mathbb{R} , with slope -1/2. The hyperbolic distance between the slope -1 and slope -1/2 lines is a constant equal to about 0.49. Call the region between these two lines a *zone*. If two paired towers have their tops in a single zone, their domains overlap. If no towers are completely truncated away between the time when the sweep line hits a point x and when there are no towers intersecting the zone above P, each equidistant curve between the two given ones hits a transmission tower at a truncated point. So in this time either the number of towers has decreased or at least 0.49 progress has been made in reducing total cost.

If we merge transmission towers when possible, a complexity based on the cost decreases sufficiently fast to give a polynomial time algorithm. The calculations in Theorem 12 implemented this geometric picture.

5. 3-MANIFOLD KNOT GENUS IS NP.

In this section we establish that 3-MANIFOLD KNOT GENUS is NP.

Proof of Theorem 2: We begin with a simplicial complex consisting of t tetrahedra whose faces are identified in pairs, and a collection of edges K in the 1-skeleton of this complex. While there are alternate formats in which a knot and a 3-manifold may be presented, all reasonable ones appear to be transformable to one another in polynomial time.

In time polynomial in t we can check that the link of each vertex is connected and has Euler characteristic two, which means that it is a sphere, and that the link of each edge is a connected curve. These are the necessary conditions to ensure that the underlying space of the complex is a 3-manifold M. Similarly we can check in time polynomial in t that the edges of K form a simple closed curve in M, and that this curve represents a trivial element of the first homology group of M with integer coefficients. We then form the second barycentric subdivision of the triangulation of M, replacing each tetrahedron by 576 tetrahedra. Removing all closed tetrahedra that meet K results in a 3-manifold M_K with a single torus boundary component, the "peripheral torus" that surrounds the knot K. The knot K bounds a surface of genus g in M if and only if there is a surface in M_K of genus g with a single boundary component that is an essential curve on ∂M_K . We restate our problem in the triangulated manifold M_K as the question of whether there exists in M_K an orientable surface of genus g with a single essential boundary component on ∂M_K . By Schubert [21], if such a surface exists, then such a surface exists among the fundamental normal surfaces in M_K . The certificate consists of an integer vector \mathbf{w} in \mathbb{Z}^{7t} giving the normal surface coordinates of this surface.

Recall that not all vectors in \mathbb{Z}^{7t} correspond to normal surfaces. It is necessary that \mathbf{w} satisfies the matching, positivity and quadrilateral conditions. These conditions can be checked for the coordinates of \mathbf{w} in time linear in $t \log W$, where W is the weight of this surface. So we can verify that there is a normal surface F with $\mathbf{w} = \mathbf{v}(F)$.

To verify that ∂F is essential on the peripheral torus, we include in the certificate a non-trivial cycle in the 1-skeleton of ∂M_K that intersects ∂F in an odd number of

points. Such a cycle can be found in the 1-skeleton of ∂M_K since curves contained in the 1-skeleton generate its first homology. The existence of such a cycle implies that ∂F is non-separating on ∂M_K , and in particular does not bound a disk on ∂M_K , ensuring that ∂F corresponds to a longitude curve parallel to K in N_K . Using the orbit counting algorithm, we can count the number of components of the normal surface F and verify that F is connected in time polynomial in $t \log N$, where N = W is the total number of points in which F meets the 1-skeleton of M_K , We can also apply Corollary 14 to verify that ∂F is connected in time bounded by a polynomial in $t \log N$. (This last step can be avoided. The number of boundary components must be odd, since ∂F represents a non-trivial element in $H_1(\partial M_K; Z_2,$ and an even number of them can be removed by joining adjacent pairs of curves with annuli on ∂M_K . This gives a surface with one boundary component and the same Euler characteristic.)

To check that F is orientable, we take the normal vector $2\mathbf{v}(F)$, that doubles each coordinate of F, and apply the orbit counting algorithm to determine if it is connected. Since M is orientable, $2\mathbf{v}(F)$ is connected if F is connected and non-orientable, and has two components if F is connected and orientable.

The Euler Characteristic of F is determined by the number of vertices, edges and faces of F, which are easily derivable from its normal coordinates. Following [11], we let t_i be the number of tetrahedra containing edge e_i and set $\epsilon_{ij} = 1$ if the edge e_i meets the j^{th} normal disk. Then the normal surface F with coordinates $\{v_j\}$ has $\chi(F) = (1/2)f_3 - \sigma(F) + wt(F)$ where $\sigma(F) = \sum v_j$ and $wt(F) = \sum_{i,j} \epsilon_{ij} x_i/t_i$. The values of ϵ_{ij}, t_i are determined by the triangulation and are independent of F. They can be computed in time polynomial in t. Since the genus of a connected orientable surface with one boundary component and Euler characteristic χ is equal to $(1 - \chi)/2$, we can determine the genus of F in time polynomial in $t \log N$.

Theorem 4 implies that the normal coordinates of this surface are at most $t2^{7t+2}$. There are 7t normal coordinates, and each represents a triangle or quadrilateral, so that the total number of intersections with the 1-skeleton satisfies $N \leq 28t^22^{7t+2}$. In particular, $\log N$ is bounded above by a polynomial in t. So the fact that F is a spanning surface for K can be verified in time polynomial in t.

6. An extended counting algorithm

In this section we develop a generalized version of the orbit counting algorithm, that counts not only the number of orbits of a collection of isometries between subintervals of an interval, but also more general quantities which are useful in applications. For example, we can use the extended algorithm to answer the following question: given a normal surface and a triangulation, how many times does each component of the surface intersect a fixed edged of the triangulation? The extended algorithm allows one to effectively compute the normal coordinates and the Euler characteristic, hence the genus, of each connected component of the surface, even when there are exponentially many components. To carry out such computations, we extend the previous analysis to pairings of weighted intervals, in which each point of the interval has associated to it a vector in \mathbb{Z}^d . We are interested in the sum of these vectors over an orbit, the orbit weights.

Consider again a pseudogroup of interval isometries acting on [1,...,N]. We will assume that there is given as input a weight function $z:[1,...,N] \to \mathbb{Z}^d$, associating to each element of [1,...,N] a vector in \mathbb{Z}^d satisfying the following condition: the

weight at successive points j and j+1 changes at most 4k times. The algorithm proceeds as before, while maintaining data on the orbit weights. We keep track of these weights by maintaining two lists of weighted subintervals. The first

$$L = \{([p_1, q_1], z_1), ..., ([p_m, q_m], z_m)\}, 1 \le p_i \le q_i < p_{i+1}$$

records the current weight values at each point in [1, N], with q_m initially equal to N. This list is updated as the algorithm proceeds. Points in the interval $[p_j, q_j]$ are assigned a constant weight $z_j \in \mathbb{Z}^d$. A second list of t subintervals

$$L' = \{([r_1, s_1], v_1), ..., ([r_t, s_t], v_t)\}, 1 \le r_i \le s_i < r_{i+1}$$

contains one point for each orbit, along with an orbit weight v_i for that orbit. Initially empty, at the algorithm's conclusion L' records the total number of orbits $\sum_{i} (s_i - r_i + 1)$, along with an orbit weight v_i assigned to each of the $(s_i - r_i + 1)$ orbits in the interval $[r_i, s_i]$.

We define an additional operation, called transferring weights by a pairing g. Suppose that $g:[a,b] \to [c,d]$ is a pairing and that [c,d] carries n different weights, given by the list

$$\{([c = r_1, s_1], v_1), ..., ([r_n, s_n = d], v_n)\}, 1 \le r_i \le s_i < r_{i+1}.$$

The weight function can be split into constant functions on n subintervals of [c,d], where $1 \le n \le d-c+1$.

The transfer operation sets the weights on [c,d] to zero and keeps the orbit weights the same by translating the weight vectors of [c,d] to smaller orbit representatives, as below:

Case 1: g is orientation preserving and b < c. Set the weights on [c, d] to zero and for each $1 \le j \le n$, add v_j to $g^{-1}([r_j, s_j])$.

Case 2: g is orientation preserving and $b \ge c$. Then $g: [a, b] \to [c, d]$ is a periodic pairing of period t = c - a. We will set the weights of points in [c, d] to zero and adjust weights in [a, c - 1] to compensate. The points in [c, d] have weights given by its intersection with intervals in L. These weights are described by

$$\{([c = r_1, s_1], v_1), ..., ([r_n, s_n = d], v_n)\}, 1 \le r_i \le s_i < r_{i+1}.$$

For each constant weight interval $([r_j, s_j], v_j)$, $1 \le j \le n$, of width w_j , add $[[\frac{w_j}{t}]]v_j$ to the weight of each point in [a, c-1], and add an additional v_j to each point in $[a+(r_j-a) \pmod{t}, a+(s_j-a) \pmod{t}]$. Finally, set the weights on [c,d] to zero.

Case 3: g is orientation reversing. We first trim g. This does not affect the orbits or the orbit weights. For a trimmed, orientation reversing pairing $g:[a,b] \to [c,d]$, set the weights on [c,d] to zero and for each $1 \le j \le n$ add v_j to $g^{-1}([r_j,s_j])$.

Lemma 15. The operation of transfer sets the weights on [c,d] to zero and preserves the orbit weights of a collection of pairings. The number of distinct weights taken by the weight function on [1,N] increases by at most four following a transfer operation.

Proof. In Cases (1) and (3), the decrease in the weight function at one point x in an orbit is exactly offset by an equal increase at the point $g^{-1}(x)$ in the same orbit.

In Case (2), each point in [a, d] is in the orbit of a unique point in [a, t - 1] = [a, c-1] under the iterates of g. The total weight of an orbit within a periodic pairing is transferred to the orbit representative in this initial subinterval by applying

powers of g. Adding weight v_j to an orbit representative in [a,c-1] of each point in $[r_j,s_j]$ while setting the weight at that point to zero, preserves the orbit weight. The resulting weights on [a,c-1] are gotten by adding appropriate multiples of v_j , with the factor being the number of orbits of a point in [a,c-1] that lie in $[r_j,s_j]$. The number of orbits under g of a point in [a,c-1] that lie in $[r_j,s_j]$ is $[[\frac{w_j}{t}]]$ or one more than this for points whose orbit hits the last $w_j \pmod{t}$ points of $[r_j,s_j]$. This latter set of points lies in the interval $[a+(r_j-a)(\bmod{t}),a+(s_j-a)(\bmod{t})]$. The formula in Case (2) follows.

Setting the weights on [c,d] to zero can cause at most two new points where a weight change occurs. The transferred weights from a constant weight interval $[r_j, s_j]$ result in a net increase of at most two pairs of successive points where the weight changes, at the preimages of its two endpoints. In Cases (1) and (3), transferred weights from an interval with non-constant weights results in an increase in the number of weight changes in the domain of g. The increase in the number of weight changes in the domain is exactly canceled by the decrease in the number of weight changes in the range of g, except possibly for two extra weight changes at the boundary points a, b of the domain. In Case (2) the same holds, but with [a, c-1] replacing the domain. In each case the number of constant weight intervals m' is increased by at most four during a transfer operation.

We now describe the modified algorithm. Again N' represents the current interval length, and we set m' to be the current number of constant weight intervals.

Weighted orbit counting algorithm: Let $\{g_i : [a_i, b_i] \to [c_i, d_i], 1 \le i \le k\}$ be a collection of pairings between subintervals of $\{1, 2, ... N\}$, and let

$$L = \{([p_1, q_1], z_1), ..., ([p_m, q_m], z_m)\}, 1 < p_i < q_i < p_{i+1} = N$$

be a list representing a collection of weights on [1, N], with the weight on $[p_j, q_j]$ equal to $z_j \in \mathbb{Z}^d$. Initialize a second weight list L' to be empty. The algorithm proceeds as before, reducing the interval size N until it reaches zero, but this time keeping track of orbit weights by maintaining the lists L, L'.

- 1. Search through the pairings and delete any pairings which are restrictions of the identity. Leave the weight lists L, L' unchanged.
- 2. Search for and contract static intervals. If the interval [r,s] is contracted, has constant weight z, and is not contained in a larger contracted interval with the same weight z, add an interval ([N'+1,N'+s-r+1],z) of width s-r+1 to the end of L', with associated weight z. Alter L by replacing [1,N'] by [1,N'-(s-r+1)], and altering each g_i by replacing any point x in a domain or range greater than s by x-(s-r+1). Replace the weight function by a new weight function w', which at points x>s satisfies w'(x)=w(x+s) and agrees with w at points x< r.
- 3. Trim all orientation reversing pairings whose domain and range overlap. Leave the weight lists L, L' unchanged.
- 4. Find a maximal pairing g_i . Recall that its range contains both N' and the range of any other pairing containing N'. If g_i is periodic, search for other periodic pairings g_j that also have range $[c_j, N']$ containing N'. If any such g_j exists, and if $t_i \leq w_j$, then perform a merger as in Lemma 10, replacing g_i and g_j by a single periodic pairing, with translation distance $GCD(t_i, t_j)$,

acting on the union of the domains and ranges of g_i and g_j . Leave the weight lists L, L' unchanged.

- 5. Find a maximal g_i . For each g_j with $j \not\models i$, if the range of g_j is contained in $[c_i, N']$, transmit g_j by g_i . Leave the weight lists L, L' unchanged.
- 6. Find the smallest value of c such that the interval [c, N'] intersects the domain or range of at most one pairing g_i , with $g_i : [a_i, b_i] \to [c_i, N']$. Transfer the weights on $[c_i, N']$ by g_i , and then truncate the pairing g_i .
- 7. If the interval size N' has decreased to zero, output the list L' and stop. Otherwise start again with Step (1).

For a \mathbb{Z}^d -valued function on [1, 2, ..., N] whose values are given by the list $L = \{([p_1, q_1], z_1), ..., ([p_m, q_m], z_m)\}, 1 \leq p_i \leq q_i < p_{i+1},$ define the total weight of L to be $\sum_{i=1}^m |z_i|$.

Theorem 16. Suppose there is an action of k pairings with \mathbb{Z}^d -valued weights, $\{g_i\}_{i=1}^k$ on $[1,\ldots,N]$, total weight bounded by D and constant weight on m subintervals. Then the weighted orbit counting algorithm outputs a list with one point for each orbit and corresponding orbit weights, and runs in time polynomial in $kmd \log D \log N$.

Proof. We will check that the running time of the algorithm is larger than that of the previous unweighted version by a factor which is a polynomial in $md \log D$.

The proof that the algorithm terminates is the same as that given for Theorem 12. There is some extra overhead involved in keeping track of weights that modifies the calculation of the running time. We indicate these additional calculations below. We now check that at each step in the algorithm the orbit weight is unchanged for any orbit remaining in L, that eliminated orbits have their orbit weights correctly recorded in L', and that k, m, d, D, N of controlled size.

As we run through the steps of the algorithm, Steps (1),(3),(4),(5) and (7) preserve the orbit structure, the weight function and the interval [1, N'], so neither of the lists L, L' are changed. The number of constant weight sub-intervals is also unchanged.

Step (2), contraction, does change the orbit structure, and also shortens [0, N']. In Step (2) the procedure adds the eliminated orbits and their weights to L'. The number of constant weight intervals is not increased, and may be decreased. Maintaining the two lists requires at most $O(md \log D)$ additional steps.

In Step (6), truncation, points are eliminated from the end of the interval [1, N']. However since we first transfer the weights of these points, the eliminated points all have weight zero, and the weight of an orbit is unaffected.

The number of steps involved in resetting the weights in L for a transfer operation is given by a polynomial in $m'd \log D \log N$. Since m' increases by at most four at each of the polynomially many steps of the algorithm, m' is bounded by a polynomial in $mk \log N$.

Combining the running time of each of the steps, whose number is given by a polynomial in $k \log N$, gives a polynomial in $kmd \log D \log N$ for the total running time.

Corollary 17. Let M be a 3-manifold with a triangulation T containing t tetrahedra and let F be a normal surface in M of total weight W. There is a procedure for counting the number of components of F and determining the topology of each component which runs in time polynomial in $t \log W$.

Proof. We begin as in Corollary 14 by assigning an integer in [1, N] to each point of intersection between the normal surface and an edge of the triangulation, where N is the total number of intersections of F with the 1-skeleton, and again associate three pairings to each face of the triangulation, one to each pair of edges in the face. The number of pairings that results is bounded above by 12t.

We next define a weight function w(x) which assigns integer weights $(z_1, z_2, \ldots, z_{7t})$ to each point in [1, N]. Initially z_i is set to zero for all i at all points $x \in [1, N]$. A tetrahedron can have as many as five distinct elementary disk types with non-zero coefficients, four triangles and one quadrilateral. If the j^{th} elementary disk type occurs, then fix one of the edges that it meets, and add 1 to the j^{th} component of the weight vector at each of the indices that the j^{th} elementary disk meets on that edge. The orbit weights are then the normal coordinates of the components of the normal surface F. Each point in the output list L' corresponds to a component of the normal surface with normal coordinates given by the corresponding weight in \mathbb{Z}^{7t} . Theorem 16 tells us that the list L' is computed in time polynomial in $kmd \log D \log N$. We now bound these constants in terms of t.

Since each edge of a tetrahedron meets at most three disk types in that tetrahedron, each edge of a tetrahedron can contribute at most six points at which the weight vector changes. Given six edges to each tetrahedron, we have $m \leq 36t$. As before we have bound for the number of pairings $k \leq 12t$ and the number of normal coordinates is given by d = 7t. The total weight bounds the normal coordinates, and with D = W we get a bound on the running time that is polynomial in $t \log W$.

Corollary 18. Let M be a 3-manifold with a triangulation T containing t tetrahedra and let F be a fundamental normal surface in M. There is a procedure for counting the number of components of F and the topology of each component which runs in time polynomial in t.

Proof. Theorem 4 gives a bound for the normal coordinates of F of $t2^{7t+2}$. Recall that there are 7t normal coordinates, and each represents a triangle or quadrilateral, so the total number of intersections with the 1-skeleton satisfies $W \leq 28t^22^{7t+2}$. In particular, $\log W$ is bounded above by a polynomial in t. Plugging this in for W in Corollary 17 we get a bound for the running time which is a polynomial in t. \square

7. The complexity of minimal spanning area

In this section we examine the complexity of the problem of determining the smallest area of a spanning surface for a curve in a 3-dimensional manifold. Such an area calculation problem seems at first to be ill suited to a complexity analysis, since it has real solutions depending on a choice of Riemannian metric.

We recast the area calculation problem into a discretized form where its complexity can be analyzed. Given a curve in a suitably discretized Reimannian 3-manifold, we ask whether it bounds a surface of area less than C, where C is an integer. To describe a metric on a 3-manifold with a finite amount of data, we restrict to piecewise flat metrics, and manifolds constructed from collections of flat tetrahedra and triangular prisms whose faces are identified by isometries. The curvature of such PL metrics is concentrated along their edges and vertices. A particular manifold in this class is described by a decomposition into tetrahedra or triangular prisms with a rational (or integer) length assigned to each edge, corresponding to edge lengths of Euclidean triangles. Prisms are allowed in order to form metrics with rational

lengths on spaces that are products. Identified 2-dimensional faces are required to be isometric. We do not require that the total angle around an edge is 2π , nor do we make any metric conditions at a vertex. Such metrics are described by a finite set of data, and can be used to approximate any Riemannian metric on a manifold. We call these objects metrized PL 3-manifolds. A curve is given as a collection of edges in the 1-skeleton of M. We will show that determining whether the smallest spanning surface for a curve in such a 3-manifold has area less than an integer C is \mathbf{NP} -hard.

Problem: MINIMAL-SPANNING-AREA

INSTANCE: A 3-dimensional metrized PL manifold M, a 1-dimensional

curve K in the 1-skeleton of M, and a natural number C.

QUESTION: Does the 1-complex bound a surface of area A < C?

The size of an instance is given by the number of bits needed to describe all the edge lengths plus $\log C$.

Theorem 19. MINIMAL-SPANNING-AREA is NP-hard.

Proof. We reduce in polynomial time an instance of the **NP**-hard problem ONE-IN-THREE SAT to an instance of MINIMAL-SPANNING-AREA. This shows that MINIMAL-SPANNING-AREA is at least as hard, up to polynomial time reduction, as ONE-IN-THREE SAT.

As a first step, we set up a 2-dimensional version of MINIMAL-SPANNING-AREA. We then construct a 3-manifold by a thickening process, with the property that a minimizing surface must remain within the 2-complex.

Given a boolean expression representing an instance of ONE-IN-THREE SAT, we construct a triangulated metrized 2-complex and an integer C. This complex contains a curve K with the property that the expression admits a satisfying assignment if and only if K bounds a surface of area less than C. This metrized complex is shown in Figure 5 for the expression $(x_1 \lor x_2 \lor x_3) \land (x_1 \lor \bar{x}_2 \lor \bar{x}_3)$.

The branched surface is similar to the one used in the proof of Theorem 1, but carries the additional structure of a metrized triangulation, whose triangles have flat metrics of prescribed edge length. The metrized triangles are constructed so that near each of the m boundary components corresponding to clauses of the boolean expression there are three triangulated disks of area close to one, one on each of the three handles coming into the punctured sphere near the boundary component. These disks are shaded in Figure 5. Each of these shaded disks is chosen to have area between 1 and 1+1/2m. The surface is constructed so that the union of all triangles in the rest of the surface has total area less than 1/2.

We saw in Theorem 1 that a spanning surface which has minimal genus goes over each of the shaded disks at most once and goes over exactly one shaded disk for each of the m clauses. It follows that such a surface has total area m < A < m+1. Furthermore, a satisfying assignment for ONE-IN-THREE SAT leads to an embedded spanning surface, with the satisfying values of the variables selecting branches of the surface, and such a spanning surface has area less than m+1. So an instance of ONE-IN-THREE SAT can be reduced to an instance of MINIMAL-SPANNING-AREA for this 2-complex.

To pass to a 3-manifold, thicken each triangle in the branched surface to a triangular prism, triangulated as in the proof of Theorem 1, and with a product metric. This produces a 3-manifold M which is a thickened up version of the 2-complex. Projection to the branched surface is area non-increasing, and area

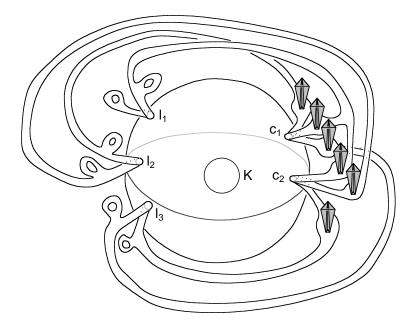


FIGURE 5. A metrized branched surface corresponding to the boolean expression $(x_1 \lor x_2 \lor x_3) \land (x_1 \lor \bar{x}_2 \lor \bar{x}_3)$. The picture is not to scale. The shaded prisms are constructed to each have area one, while the rest of the surface has total area less than 1/2.

decreasing for a surface with boundary on the branched surface but not contained in it. Therefore a least area surface spanning K must lie on the branched surface. A closed manifold DM with a piecewise-smooth metric can be obtained by a doubling construction as in Theorem 1. The doubling involution is an isometry, so that reflecting a surface meeting $DM \setminus M$ into M does not increase area. It follows that the embedded spanning surface on the branched surface is a least area surface in DM.

8. Open questions

Among many unresolved questions are:

- 1. Does determining knot genus remain **NP**-hard if we restrict to knots in the 3-sphere?
- 2. Is determining the genus of a knot in a 3-manifold \mathbf{NP} ? This amounts to showing that finding a lower bound to the knot's genus is an \mathbf{NP} problem, in contrast to the upper bound we have investigated. Recall that the genus of a knot is the least possible genus of all spanning surfaces. We have shown that certifying that the genus is at most g is \mathbf{NP} , but left open the possibility that the genus may be smaller than g. If the answer to this question is yes, then we can certify that a non-trivial knot has positive genus, and it would follow that UNKNOTTING is both \mathbf{NP} and \mathbf{coNP} .

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