

## 4.4

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# $\kappa$ -Compatible Tessellations\*

Philippe P. Pébay and David Thompson

Sandia National Laboratories  
P.O. Box 969, Livermore CA 94551, U.S.A.  
{pppebay,dcthomp}@sandia.gov

**Summary.** The vast majority of visualization algorithms for finite element (FE) simulations assume that linear constitutive relationships are used to interpolate values over an element, because the polynomial order of the FE basis functions used in practice has traditionally been low – linear or quadratic. However, higher order FE solvers, which become increasingly popular, pose a significant challenge to visualization systems as the assumptions of the visualization algorithms are violated by higher order solutions. This paper presents a method for adapting linear visualization algorithms to higher order data through a careful examination of a linear algorithm’s properties and the assumptions it makes. This method subdivides higher order finite elements into regions where these assumptions hold ( $\kappa$ -compatibility). Because it is arguably one of the most useful visualization tools, isosurfacing is used as an example to illustrate our methodology.

## 1 Introduction

People have been approximating solutions to partial differential equations (PDEs) ever since PDEs were conceived. Much more recently, a class of techniques known as *hp*-adaptive methods have been developed in an effort to converge to a solution faster than previously possible, or to provide more accurate approximations than traditional finite element simulation within the same amount of computational time. These new techniques can increase both the hierarchical (*h*) and polynomial (*p*) levels of detail – or degrees of freedom – during a simulation.

Once these solution approximations have been computed, they must be characterized in some way so that humans can understand and use the results. This paper develops a technique for partitioning higher-order cells in order to

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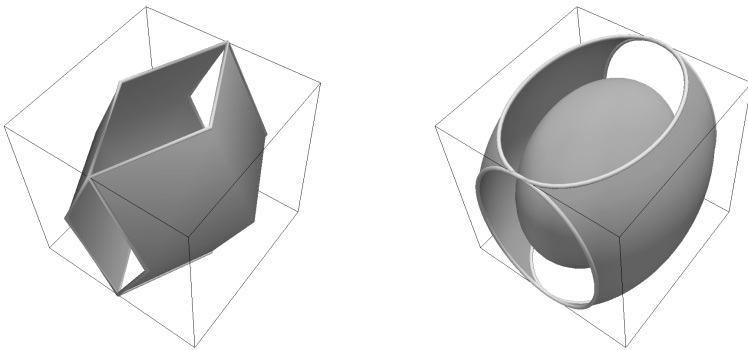
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characterize the behavior of their geometric and scalar field curvatures during post-processing. We say that such partitions are  $\kappa$ -compatible. While a past paper [8] has presented our software framework for creating these partitions, this paper presents a full description of the algorithm and a rigorous proof of the conditions under which it will work and terminate.

Currently, visualization techniques for quadratic and higher order FE solutions are very limited in scope and/or cannot guarantee that all topological features are captured [1, 3, 7, 8, 6]. These techniques are also limited in their applicability to a subset of visualization techniques. Moreover, although some production-level tools currently offer support for quadratic elements, they do not always do so correctly (*cf.* [8] for a discussion); in the case of isocontouring, consider for example the following scalar field:

$$\begin{aligned} \Phi_1 : [-1, 1]^3 &\longrightarrow \mathbb{R} \\ (x, y, z)^T &\mapsto x^2 + y^2 + 2z^2, \end{aligned}$$

interpolated over a single Q2 Lagrange element (hexahedron with degree 2 Lagrange tensor-product interpolation over 27 nodes), with linear geometry, where one is interested in the isocontours  $\Phi_1^{-1}(1)$  and  $\Phi_1^{-1}(2)$ .



**Fig. 1.** Isocontours of  $\Phi_1$  for the isovalues 1 (cyan) and 2 (green): linear isocontouring approach (left), and our new topology-based approach (right).

As shown in Figure 1, left, the linear isocontouring approach (implemented here in ParaView) completely misses  $\Phi_1^{-1}(1)$ , because it is entirely contained within the cell. Meanwhile, the topology of  $\Phi_1^{-1}(2)$  is correct, but its geometry is poorly captured. On the other hand, our new method (Figure 1, right) captures the correct topologies of both isocontours, and provides a much better geometric approximation of  $\Phi_1^{-1}(2)$  than linear isocontouring does.

The lack of tools that are applicable to most visualizations of higher order element simulations, and that can guarantee correctness of the results, prevents analysts from exploiting such simulations. In this paper, we propose a solution to this problem.

### 1.1 Higher Order Finite Elements

Here we briefly review the FE method to develop notation used throughout the paper. Recall that the FE method approximates the solution,  $f : \Omega \mapsto \mathbb{R}$ , of some PDE as a set of piecewise functions over the problem domain,  $\Omega \subset \mathbb{R}^d$ . Although  $\Omega$  may be any general  $d$ -dimensional domain, we'll assume it is 3-dimensional. The fact that we have a piecewise approximant divides  $\Omega$  into subdomains  $\Omega_e \subseteq \Omega$ ,  $e \in E_\Omega$  that form a closed cover of  $\Omega$ . Each  $\Omega_e$  is itself a closed set that is parametrized with a map (usually polynomial in form) from parametric coordinates  $\mathbf{r} = (r, s, t) \in R \subset \mathbb{R}^3$  to geometric coordinates  $\mathbf{x} = (x, y, z) \in \Omega_e$ :

$$\Xi_e(\mathbf{r}) = B_{0,0,0} + B_{1,0,0}r + B_{0,1,0}s + B_{0,0,1}t + B_{1,1,1}rst + \dots$$

where  $B_{i,j,k} \in \mathbb{R}^3$  such that  $\Xi_e$  is invertible for all points of  $\Omega_e$ . Therefore, the approximate solution may be written in terms of parametric coordinates (as  $\Phi_e$ ) or in terms of geometric coordinates:  $f_e(\mathbf{x}) = \Phi_e \circ \Xi_e^{-1}(\mathbf{x})$ , where

$$\begin{aligned} \Phi_e : R \subset \mathbb{R}^3 &\longrightarrow \mathbb{R} \\ (r, s, t)^T &\mapsto \Phi_e(r, s, t) \end{aligned}$$

is the approximating function over  $\Omega_e$  expressed in terms of parametric coordinates. Furthermore, we require that each  $\Phi_e$  is polynomial in the parameters:

$$\Phi_e(\mathbf{r}) = A_{0,0,0} + A_{1,0,0}r + A_{0,1,0}s + A_{0,0,1}t + A_{1,1,1}rst + \dots$$

These coefficients,  $A_{i,j,k} \in \mathbb{R}$ , are known as *degrees of freedom* (DOFs). Each one corresponds to a particular modal shape, and sets of modal shapes can be grouped together into nodes by the regions of parameter-space over which they have an effect (a given corner, edge, face, or the interior volume).

A global approximant  $f(\mathbf{x})$  can then be constructed from the piecewise elemental approximants. This leaves only the matter of what to do where  $\Omega_e$  and  $\Omega_{j,j \neq e}$  intersect. Usually, these subdomains intersect over  $(d - 1)$ -dimensional or lower regions (2-dimensional faces, 1-dimensional edges, and “0”-dimensional vertices in our case). In these regions,  $\Phi$  is not well-defined since  $\Phi_e$  and  $\Phi_j$  may disagree. Usually, the FE method constrains  $\Phi_e$  and  $\Phi_j$  to be identical in these regions, however some methods such as the discontinuous Galerkin method do not require this and subsequently have no valid approximant in these regions.

For a given decomposition of  $\Omega$ , the FE method may not converge to the correct (or indeed, any) solution. When  $\Phi_e$  and  $\Xi_e$  are trilinear polynomials for all  $e$ , a technique called  $h$ -adaptation is often used to force convergence and/or increase solution accuracy. In this technique, some subdomains  $E \subseteq E_\Omega$  are replaced with a finer subdivision  $E'$  such that  $\bigcup_{e \in E} \Omega_e = \bigcup_{e \in E'} \Omega_e$  but  $|E'| > |E|^2$ . Similarly,  $p$ -adaptation is the technique of increasing the

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<sup>2</sup>In temporal simulations, we do allow  $|E'| < |E|$  in regions where  $\Omega$  has been adapted to some time-transient phenomenon.

order of polynomials  $\Phi_e$  and/or  $\Xi_e$  rather than the number of finite elements.  $hp$ -adaptation is then some combination of  $h$ - and  $p$ -adaptation over  $\Omega$ . In the end, the FE method provides an approximation to  $f = \Phi \circ \Xi^{-1}$  by solving a collection of equations for coefficients ( $A$  and  $B$  in the examples above). Our task is then to characterize the maps  $\Phi_e$  and  $\Xi_e$  in a way that aids human understanding of the solution.

The rest of this paper presents a method for partitioning higher-order finite elements into regions where visualization assumptions are satisfied.

## 2 Partitioning Finite Elements

Using these notation, we now examine what requirements must be satisfied by  $\kappa$ -compatible tessellations. A common assumption made by linear visualization algorithms is that critical points (points where all partial derivatives of  $f_e$  vanish) may only occur at vertices. This one assumption can show up in many different ways. Algorithms that iterate over an element's corner vertices to identify extrema (*e.g.*, thresholding) all make this assumption. Other examples include (but are not limited to) isosurfacing, cutting, and clipping.

This paper is concerned with tessellating finite elements into regions where the critical-points-at-vertices assumption holds as part of an overall strategy to adapt existing techniques to work with higher order elements. Because we use it as a running example of an important application driving the development of  $\kappa$ -compatible tessellation, we will focus on the prerequisites of the linear tetrahedral isosurfacing algorithm, keeping in mind that other visualization techniques result in the same set of constraints. The linear tetrahedral isosurfacing algorithm assumes:

- (C1<sup>0</sup>) each tetrahedron edge intersects a particular isocontour at most once,
- (C2<sup>0</sup>) no isocontour intersects a tetrahedron face without intersecting at least two edges of the face,
- (C3<sup>0</sup>) no isocontour is completely contained within a single tetrahedron, and
- (C4<sup>0</sup>) the map from parametric to geometric coordinates must be bijective.

*Remark 1.* Note that (C4<sup>0</sup>) only regards  $\Xi_e$  and can be restated as:

$$(C4) \forall \mathbf{x} \in \Omega_e, \exists! \mathbf{r} \in R \text{ such that } \Xi_e(\mathbf{r}) = \mathbf{x}.$$

In this paper, it is satisfied by hypothesis as we focus on the  $\Phi_e$  map. However, in the context of higher order element mesh generation and modification, our methodology can be applied to  $\Xi_e$  to verify the correctness of each element.

Let's examine how changing to a higher-order interpolant affects these assumptions. For instance, the  $\Phi_1^{-1}(1)$  isocontour in the example of §1 violates (C3<sup>0</sup>), which explains why it is entirely missed by linear isocontouring.

We now translate the criteria into requirements on  $\Phi_e$ , that are slightly stronger for reasons that are discussed later on.

**Proposition 1.**  $(C1^0)$ ,  $(C2^0)$  and  $(C3^0)$  are respectively implied by:

(C1) for each edge  $E$  of  $R$ , with direction vector  $(a_x, a_y, a_z)$ ,

$$a_x \frac{\partial \Phi_e}{\partial r} + a_y \frac{\partial \Phi_e}{\partial s} + a_z \frac{\partial \Phi_e}{\partial t} \neq 0 \quad \text{over the interior of } E.$$

(C2) for each face  $F$  of  $R$ , with vector basis  $((a_x, a_y, a_z), (b_x, b_y, b_z))$ ,

$$\begin{cases} a_x \frac{\partial \Phi_e}{\partial r} + a_y \frac{\partial \Phi_e}{\partial s} + a_z \frac{\partial \Phi_e}{\partial t} \neq 0 \\ b_x \frac{\partial \Phi_e}{\partial r} + b_y \frac{\partial \Phi_e}{\partial s} + b_z \frac{\partial \Phi_e}{\partial t} \neq 0 \end{cases} \quad \text{over the interior of } F.$$

(C3)  $\nabla \Phi_e \neq 0$  over the interior of  $R$ .

*Proof.* By definition,  $(C1^0)$ ,  $(C2^0)$  and  $(C3^0)$  are equivalently stated as:

(C1<sup>0</sup>) No restriction of  $\Phi_e$  to an element edge has extrema interior to the edge unless the restriction is constant on the edge.

(C2<sup>0</sup>) No restriction of  $\Phi_e$  to an element face has extrema interior to the face unless the restriction is constant on the face.

(C3<sup>0</sup>)  $\Phi_e$  has no extrema inside the interior of the element unless the interpolant is constant over the entire element.

Being a polynomial function,  $\Phi_e$  is  $C^1$ , and so are its restrictions to the edges and faces of  $R$ . Let  $E$  be an edge of  $R$  that passes through the point  $(p_x, p_y, p_z)$ , with direction vector  $(a_x, a_y, a_z)$ . This edge can be parametrized as follows:

$$\begin{aligned} \eta : I \subset \mathbb{R} &\longrightarrow \mathbb{R}^3 \\ t &\longmapsto (a_x t + p_x, a_y t + p_y, a_z t + p_z)^T, \end{aligned}$$

from which we obtain the restriction of  $\Phi_e$  to  $E$ ,  $\Phi_{e|E} = \Phi_e \circ \eta : I \longrightarrow \mathbb{R}$ . Its derivative  $d\Phi_{e|E} \in \mathcal{L}(\mathbb{R})$  at any arbitrary point  $t_0 \in I$  is thus

$$d\Phi_{e|E}(t_0) = d\Phi_e(\eta(t_0)) \circ d\eta(t_0) = \left( \frac{\partial \Phi_e}{\partial r}, \frac{\partial \Phi_e}{\partial s}, \frac{\partial \Phi_e}{\partial t} \right) \Big|_{\eta(t_0)} (a_x, a_y, a_z)^T dt. \quad (1)$$

As  $\Phi_{e|E}$  is  $C^1$  over  $I$ , a necessary condition for  $\Phi_{e|E}$  to have an extremum in  $t_0$ , interior to  $I$ , is that  $d\Phi_{e|E}(t_0) = 0$ , i.e.,

$$a_x \frac{\partial \Phi_e}{\partial r}(t_0) + a_y \frac{\partial \Phi_e}{\partial s}(t_0) + a_z \frac{\partial \Phi_e}{\partial t}(t_0) = 0.$$

Similarly, let  $F$  be a face of  $R$  that passes through the point  $(p_x, p_y, p_z)$  and has basis  $((a_x, a_y, a_z), (b_x, b_y, b_z))$ . This face can be parametrized with two variables as follows:

$$\begin{aligned} \eta : U \subset \mathbb{R}^2 &\longrightarrow \mathbb{R}^3 \\ (u, v)^T &\longmapsto (a_x u + b_x v + p_x, a_y u + b_y v + p_y, a_z u + b_z v + p_z)^T. \end{aligned}$$

The derivative  $d\Phi_{e|F} \in \mathcal{L}(\mathbb{R}^2, \mathbb{R})$  of  $\Phi_{e|F} = \Phi_e \circ \eta : U \rightarrow \mathbb{R}$  is

$$d\Phi_{e|F}(u_0, v_0) = d\Phi_e(\eta(u_0, v_0)) \circ d\eta(u_0, v_0) \tag{2}$$

$$= \left( \frac{\partial \Phi_e}{\partial r}, \frac{\partial \Phi_e}{\partial s}, \frac{\partial \Phi_e}{\partial t} \right) \Big|_{\eta(u_0, v_0)} \begin{pmatrix} a_x & b_x \\ a_y & b_y \\ a_z & b_z \end{pmatrix} \begin{pmatrix} du \\ dv \end{pmatrix}. \tag{3}$$

As  $\Phi_{e|F}$  is  $C^1$  over  $U$ , a necessary condition for  $\Phi_{e|F}$  to have an extremum in  $(u_0, v_0)$ , interior to  $U$ , is that  $d\Phi_{e|F}(u_0, v_0) = 0$ , which may be restated as

$$\begin{aligned} a_x \frac{\partial \Phi_e}{\partial r}(\eta(u_0, v_0)) + a_y \frac{\partial \Phi_e}{\partial s}(\eta(u_0, v_0)) + a_z \frac{\partial \Phi_e}{\partial t}(\eta(u_0, v_0)) &= 0 \\ b_x \frac{\partial \Phi_e}{\partial r}(\eta(u_0, v_0)) + b_y \frac{\partial \Phi_e}{\partial s}(\eta(u_0, v_0)) + b_z \frac{\partial \Phi_e}{\partial t}(\eta(u_0, v_0)) &= 0. \end{aligned}$$

Finally, as  $\Phi_e$  is  $C^1$  over  $R$ , a necessary condition for  $\Phi_e$  to have extremum in  $(r_0, s_0, t_0)$ , interior to  $R$ , is that  $d\Phi_e(r_0, s_0, t_0) = 0$ , i.e.,  $\nabla \Phi_e(r_0, s_0, t_0) = 0$ .  $\square$

Note that the proof mainly relies on the fact that, for a  $C^1$  function over an open domain, an extremum is a critical point. However, the converse is not true and thus each  $(Ci)$  is stronger than the corresponding  $(Ci^0)$ . Therefore, using  $(Ci)$  rather than  $(Ci^0)$  can yield non-extremal critical points – just consider  $r \mapsto r^3$  in 0. In order to eliminate such “false positives”, we would need to evaluate second- and higher-order derivatives, which would incur further computational costs, and even this would not always suffice as degenerate cases may occur. Therefore, rather than degrading computational performance, the trade-off we make is to accept the stronger  $(Ci)$  conditions and extra points that are not required by the  $(Ci^0)$  conditions. This may even be beneficial in terms of geometric approximation, as non-extremal critical points can be the locus of important geometric features (e.g., saddle points). In short,  $(Ci)$  conditions mean that all of the differences between the linear and higher order isocontouring implementations can be attributed to critical points of  $\Phi_e$ .

### 2.1 Creating the Partition

In §2, we presented the requirements a tetrahedral element must meet for isocontouring algorithm to work. As we noted earlier, higher order elements that have non-simplicial domains (such as hexahedra, pyramids, etc.) will have to be decomposed into tetrahedra  $\mathcal{T}$ . However, these tetrahedra must additionally meet the requirements (C1) to (C4). As explained in Remark 1, (C4) is satisfied by hypothesis; therefore the partition is subdivided until it meets criteria (C1), (C2), and (C3). Because not only a single scalar field, but a set  $\kappa$  of such fields may be of interest, these criteria must be satisfied for all fields in  $\kappa$ . Once this is achieved, the final partition is said to be  $\kappa$ -compatible. For the sake of legibility, we only discuss here the case where  $\kappa = \{\Phi\}$ , but the method remains the same when  $\kappa$  is not a singleton.

From now onwards, **it is assumed that all critical points are isolated.** This requirement is necessary so that the set of all critical points is finite, since a polynomial function can only have a finite number of isolated critical points; its implications are discussed at the end of this section. In this context, the general scheme of our method applied to the input parameters  $M$  (initial mesh) and  $\Phi$  (field interpolated over  $M$ ) is

PARTITION-MESH( $M, \Phi$ )

```

1   $C \leftarrow$  DOF-CRITICALITIES( $M, \Phi$ )
2   $(T_0, S) \leftarrow$  TRIANGULATE-BOUNDARIES( $M, C$ )
3  CORRECT-TRIANGLE-TOPOLOGY( $M, \Phi, S, T_0$ )
4   $(T_1, S) \leftarrow$  TETRAHEDRALIZE-INTERIOR( $M, T_0, C$ )
5  CORRECT-TETRAHEDRAL-TOPOLOGY( $M, \Phi, S, T_1$ )
6  return  $T_1$ 

```

The output of the scheme is a tetrahedral subdivision  $T_1$  of  $M$ . We now discuss each step in details, and theoretically establish the validity of the approach.

DOF-CRITICALITIES

This algorithm locates the critical points of  $\Phi$  inside each element and of the restrictions of  $\Phi$  to all element faces and edges. It takes  $M$  and  $\Phi$  as inputs and yields a set  $C$  of critical points.  $\text{BDY}^1(R)$  and  $\text{BDY}^2(R)$  respectively denote the set of 1- and 2-dimensional boundaries of the parametric domain  $R$ .

DOF-CRITICALITIES( $M$ )

```

1  for  $e \leftarrow |M|$ 
2      do Find critical points of  $\Phi_e$  in  $R$ 
3          Store critical points indexed by volumetric DOF node.
4          for  $f_i^2 \in \text{BDY}^2(R)$ 
5              do if  $\nabla \Phi_{e|f_i^2} = 0$  not marked,
6                  then Find critical points of  $\Phi_{e|f_i^2}$ 
7                      Store critical points of  $\Phi_{e|f_i^2}$  in  $C_{f_i^2}$ 
8                      Mark  $\Phi_{e|f_i^2}$  as done
9          for  $f_i^1 \in \text{BDY}^1(R)$ 
10             do if  $\nabla \Phi_{e|f_i^1} = 0$  not marked,
11                 then Find critical points of  $\Phi_{e|f_i^1}$ 
12                     Store critical points of  $\Phi_{e|f_i^1}$  in  $C_{f_i^1}$ 
13                     Mark  $\Phi_{e|f_i^1}$  as done
14  return  $C = (\cup_i C_{f_i^1}) \cup (\cup_i C_{f_i^2})$ 

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Note that, because finding critical points is a time-consuming process, we do not wish to process shared edges or faces multiple times. This extra work is avoided by storing critical points indexed by the DOF with which they are associated – critical points on a face are stored with the index used to

retrieve the coefficients for that face’s degrees of freedom, and likewise for edges. Therefore, DOF-CRITICALITIES operates on the mesh a whole, and not independently on each element. The issue of how to actually find the critical points is a complex and challenging problem of its own. We have not specifically researched this issue, and we handle it as follows:

- for edge critical points, where the problem amounts to finding all roots of a polynomial in a bounded interval, we have implemented exact solvers for up to quartic equations (and hence, quintic interpolants), and a Lin-Bairstow solver for higher order equations;
- for face and body critical points, where the problem amounts to finding all roots of a polynomial system within a bounded domain, we solve exactly if the system is linear, and otherwise make use of the PSS package [4, 5]. However, we think that this aspect deserves much further investigation.

If one assumes that the polynomial system solver always terminates in finite time, then DOF-CRITICALITIES does as well. As mentioned earlier, because restrictions of the field to edges and faces are marked as they are done, each is processed only once even when it is shared by several elements.

*Remark 2.* The methodology we present requires the ability to detect all critical points on arbitrary line segments and triangular faces in the domain of an element. Most polynomial system solvers require a power-basis representation of a system to be solved and that is not usually how finite elements are represented. Given that we wish to perform this change of basis as infrequently as possible, it behooves us to find a way to derive the restriction of  $\Phi_e$  to a line or face from the full representation of  $\Phi_e$ .

## TRIANGULATE-BOUNDARIES

Once the critical points have been located, the second step of our scheme consists in triangulating the two-dimensional boundaries of all elements. This ensures that all volumetric elements that reference a particular face use the same triangulation – otherwise our model could have cracks along element boundaries<sup>3</sup>. In order to satisfy (C2) on a given element  $e$ , a restriction of  $\Phi_e$  to a face of the tessellation of  $e$  is not permitted to have a critical point; therefore, we “eliminate” the critical points of the restriction of  $\Phi_e$  to the faces of  $e$  by inserting them in the list of points to be triangulated. The algorithm TRIANGULATE-BOUNDARIES thus takes the mesh  $M$  and its related set of critical points  $C$  as inputs, and returns a triangulation  $T_0$  of the set of faces in  $M$ . In this algorithm, the method FACE-CENTER takes a face as its input and returns its parametric center, and STAR<sup>2</sup>( $c, Q_1, \dots, Q_n$ ) creates a

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<sup>3</sup>Discontinuous Galerkin elements can be accommodated by using different indices (as opposed to a shared index  $i$ ) for edges and faces of adjoining elements. Cracks would occur, but they would be faithful representations of the interpolant discontinuity.



triangulation composed of triangles  $cQ_1Q_2, \dots, cQ_nQ_1$  (with the requirement that  $c$  is contained in the interior of the convex hull of  $\{Q_1, \dots, Q_n\}$ ).

TRIANGULATE-BOUNDARIES( $M, C$ )

```

1  for  $f_i^2 \leftarrow$  each 2-boundary of every 3-D finite element
2      do if  $|C_{f_i^2}| > 0$ 
3          then  $c \leftarrow C_{f_i^2,0}$ 
4          else  $c \leftarrow$  FACE-CENTER(BDY $_i^2(\mathbb{R})$ )
5           $T_i \leftarrow \emptyset$ 
6           $Q \leftarrow$  corner points of face  $i \cup$  isolated critical points
           of all bounding edges of face  $f_i^2$ , ordered in a
           counterclockwise loop around  $f_i^2$ .
7          for  $j \in \{0, \dots, |Q| - 1\}$ 
8              do Insert STAR $^2(c, Q_j, Q_{(j+1) \bmod |Q|})$  into  $T_i$ 
9           $C_{f_i^2}' \leftarrow \{C_{f_i^2} \setminus C_{i,0}\}$ 
10         for  $c \in C_{f_i^2}'$ 
11             do Find  $t \in T_i$  such that  $c \in t$ 
12                 Remove  $c$  from  $C_{f_i^2}'$ 
13                 Remove  $t$  from  $T_i$ 
14                 Subdivide  $t$  into 2 or 3 triangles  $t_k$ 
15                 Insert  $t_k$  into  $T_i$ 
16  return  $T_0 = \cup_i T_i$ 

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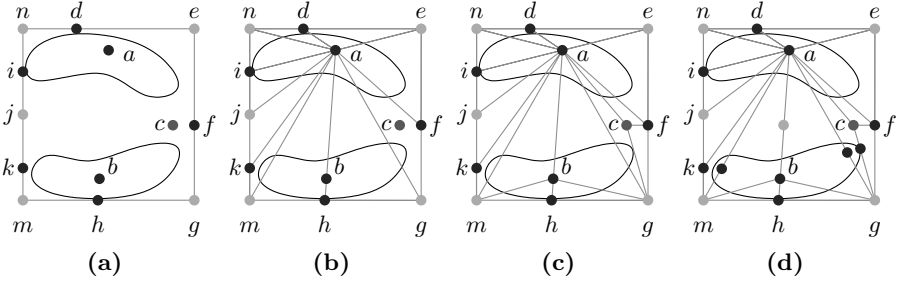
All sets involved in TRIANGULATE-BOUNDARIES are finite, and no recursion is involved. Therefore, this procedure terminates in finite time. In addition,

**Proposition 2.** *Upon completion of TRIANGULATE-BOUNDARIES, (C1) is satisfied on the edges of the triangulation  $T_0$  that either belong to  $M$  or are subdivisions of edges or faces of  $M$ , and (C2) is satisfied across  $T_0$ .*

*Proof.* By construction, TRIANGULATE-BOUNDARIES inserts in  $T_0$  all critical points of restrictions of  $\Phi$  to edges of  $M$ . Since this process cannot create novel critical points on these edges, the first part of the proposition ensues. Compliance with (C2) across  $T_0$  is ensured because all critical points of restrictions of  $\Phi$  to faces of  $M$  have been inserted as vertices of  $T_0$ , and no new such critical point may have been created.  $\square$

Note that that there is no guarantee that (C1) is satisfied across  $T_0$  upon completion of TRIANGULATE-BOUNDARIES, as shown in the following example:

*Example 1.* Consider a face  $f_i$ , illustrated in Fig. 2(a), such that the restriction of the field to the interior of  $f_i$  has 3 critical points (denoted  $a, b$ , and  $c$ ), and each of the restrictions of the field to the edges of  $f_i$  has at least one critical point (denoted  $d, f, h, i, j$ , and  $k$ ). The triangulation displayed in Fig. 2(b) is obtained once all STAR $^2$  procedures have performed by taking  $a$  as the first internal critical point to be inserted. The final tessellation, shown in Fig. 2(c), has eliminated the remaining interior critical points  $b$  and  $c$  by making them



**Fig. 2.** An example face  $f_i^2$  with critical points shown as blue dots (maxima), red dots (saddles), and green dots (minima). (a) The input to TRIANGULATE-BOUNDARIES. (b) Resulting triangulation of  $f_i^2$  after all STAR<sup>2</sup> procedures of TRIANGULATE-BOUNDARIES have been performed. (c) The triangulation at the completion of TRIANGULATE-BOUNDARIES. (d) The new critical points introduced by the first stage of CORRECT-TRIANGLE-TOPOLOGY.

nodes of the triangulation; however, as illustrated in Fig. 2(d), new critical points have appeared, on the restrictions of  $\Phi_e$  to edges  $ab$ ,  $ag$ ,  $am$ , and  $cg$ .

*Remark 3.* Line 14 of TRIANGULATE-BOUNDARIES allows for 2 different ways to subdivide a triangle, depending on whether the face critical point lies within or on the boundary of the triangle; for instance, in Fig. 2(c), triangles  $ahm$  and  $afg$  are split in, respectively, 2 and 3 triangles. In practice, to avoid unnecessary creation of quasi-degenerate triangles, the implementation of TRIANGULATE-BOUNDARIES uses a predefined threshold (that can be related to the distance of the critical point to the closest triangle edge, or to a triangle quality estimate of the subdivided triangles) below which a critical point is moved to the appropriate edge; for instance, in the example illustrated in Fig. 2(c), point  $b$  is considered as belonging to  $ah$ , even if it slightly off.

CORRECT-TRIANGLE-TOPOLOGY

This algorithm searches for critical points in the restriction of  $\Phi_e$  to each unmarked edge of  $T_0$ . When points are found, they are inserted into  $T_0$ , and iteratively repeats the procedure until (C1) and (C2) are satisfied throughout the entire triangulation of the faces of  $M$ . Fig. 3(b) shows this procedure applied to Example 1.

```

CORRECT-TRIANGLE-TOPOLOGY( $M, \Phi, S, T_0$ )
1  while  $S$  not empty
2      do Pop  $t$  from  $S$ 
3           $C \leftarrow \emptyset$ 
4          for  $e \leftarrow$  marked edges of  $t$ 
5              do Insert critical points of  $e$  into  $C$ 
6          for  $c \leftarrow C$ 
7              do Find  $t \in T_1$  s. t.  $c \in t$ 
8                  Remove  $t$  from  $T_1$  and  $S$ 
9                   $U \leftarrow \text{STAR}^2(c, t)$ 
10             for  $t' \leftarrow U$ 
11                 do if MARK-TRIANGLE( $t'$ )
12                     then Push  $t'$  onto  $S$ 
13                     Insert  $t'$  into  $T_0$ 

```

**Proposition 3.** *If, for all  $e$  in  $M$ , all critical points of the restriction of  $\Phi_e$  to any face of  $e$  are isolated, then CORRECT-TRIANGLE-TOPOLOGY terminates. In addition, upon termination, (C1) and (C2) are satisfied across  $T_0$ .*

*Proof.* For brevity, the proof is not provided here: it can readily be obtained by reducing the proof of Proposition 5 to the 2-dimensional case.  $\square$

*Remark 4.* Because there is no need to refine below a certain size for visualization purposes, our implementation uses a tolerance  $\varepsilon \in \mathbb{R}_+$  so that, in line 5, a new critical point  $p$  of the restriction of  $\Phi$  to a marked edge is inserted only if there is no  $p'$  in  $C$  such that  $|p - p'| < \varepsilon$ . Therefore, the procedure may terminate before (C1) and (C2) are fully satisfied, but are satisfied up to  $\varepsilon$ .

#### TETRAHEDRALIZE-INTERIOR

Each element interior is now tetrahedralized, and although we treat the whole mesh, there is here no issue of inter-element consistency, as this part of the scheme only regards element interiors. As in TRIANGULATE-BOUNDARIES, we “eliminate” critical points of  $\Phi_e$  that are interior to  $e$  by adding them to the set of points to be tetrahedralized. Therefore, the tetrahedralization of each element  $e$  is constrained by the triangulations of the faces of  $e$  that result from CORRECT-TRIANGLE-TOPOLOGY, and by the critical points of  $\Phi_e$  that are interior to  $e$ . Additionally, when the finite element is starred into a set of tetrahedra, we know that the triangular base of each tetrahedron and its 3 bounding edges will not have any critical points since those have already been identified and inserted into the triangulation of the two-dimensional boundary of the element. However, the remaining 3 faces and 3 edges must be marked because  $\Phi_e$  restricted to their domain may contain critical points. This is accomplished by MARK-TETRAHEDRON, which sets a bit code for each edge and face not on the base of the given tetrahedron (which must be properly oriented when passed to the subroutine). The algorithm is then as follows:

TETRAHEDRALIZE-INTERIOR( $M, T_0, C$ )

```

1   $S \leftarrow \emptyset$ 
2   $T_1 \leftarrow \emptyset$ 
3  for  $e \leftarrow |M|$ 
4      do Let  $T \subseteq T_0$  be all triangles on  $\text{BDY}^2(R)$ 
5          if  $|C_e| > 0$ 
6              then  $c \leftarrow C_{e,0}$ 
7              else  $c \leftarrow \text{ELEMENT-CENTER}(R)$ 
8           $V \leftarrow \text{STAR}^3(c, T)$ 
9          for  $t \leftarrow V$ 
10             do if MARK-TETRAHEDRON( $t$ )
11                 then Push  $t$  onto  $S$ 
12         for  $c \in \{C_e \setminus C_{e,0}\}$ 
13             do Find  $t \in V$  s. t.  $c \in t$ 
14                 Remove  $t$  from  $V$  and  $S$ 
15                  $U \leftarrow \text{STAR}^3(c, t)$ 
16                 for  $t' \leftarrow U$ 
17                     do if MARK-TETRAHEDRON( $t'$ )
18                         then Push  $t'$  onto  $S$ 
19                     Insert  $t'$  into  $V$ 
20             Insert  $V$  into  $T_1$ 
21 Return  $(T_1, S)$ 

```

All sets involved in TETRAHEDRALIZE-INTERIOR are finite, and no recursion is involved. Therefore, this procedure terminates in finite time. In addition,

**Proposition 4.** (C3) is satisfied across the tetrahedralization  $T_1$  upon completion of TETRAHEDRALIZE-INTERIOR. Moreover, any sub-tetrahedralization of  $T_1$  satisfies (C3) as well.

*Proof.* Let  $p$  be a critical point of the field  $\Phi$ ; then, 2 cases may occur:

1.  $p$  is contained in the interior of an element  $e \in M$ . In this case,  $p$  belongs to  $C_e$  (and to no other  $C_{e'}$ ) and thus, thanks to  $\text{STAR}^3$ ,  $p$  is a tetrahedron vertex in  $T_1$ .
2.  $p$  is contained on the boundary of an element  $e \in M$ . In this case, it is also a lower-dimensional critical point, *i.e.*, a critical point for the restriction of  $\Phi_e$  to one of its faces or edges, because the fact  $d\Phi_e$  vanishes in  $p$  ensures that the left hand side of (1) (if  $p$  is on an edge) or (2) (if  $p$  is on a face) vanishes as well. Therefore,  $p$  belongs to  $C_{f_i}$  for an edge  $f_i^1$  or a face  $f_i^2$  and hence has been made a triangle vertex of  $T_0$  by TRIANGULATE-BOUNDARIES. Since all triangles of  $T_0$  become tetrahedral faces in  $T_1$ , then  $p$  is a tetrahedron vertex in  $T_1$ .

In both cases, upon completion of TETRAHEDRALIZE-INTERIOR,  $p$  is a tetrahedron vertex in  $T_1$ . Therefore, for all  $t \in T_1$ ,  $p$  is not a critical point of  $\Phi_t$  interior to  $t$ . As this is true for any critical point  $p$  of the field  $\Phi$ , it follows

that  $T_1$  satisfies (C3). Finally, as (C3) is satisfied on any tetrahedron  $t \in T_1$ , it is also satisfied on any tetrahedral subdivision of  $t$ : indeed, if one could find a tetrahedron  $t' \subset t$  and a critical point  $p$  of  $\Phi$  contained in the interior of  $t'$ , then  $p$  would also be in the interior of  $t$ , which would contradict the hypothesis; *ad absurdum*, the result ensues.  $\square$

#### CORRECT-TETRAHEDRAL-TOPOLOGY

At this point, a brief summary of what has been obtained through the 3 first stages of our scheme will most likely be useful to the reader. A tetrahedralization  $T_1$  of the initial mesh  $M$  has been obtained, such that

- (C3), and (C4) by hypothesis, are satisfied;
- (C1) and (C2) are satisfied on all edges and faces of  $T_1$  that either belong to  $M$  or are subdivisions of edges or faces of  $M$ .

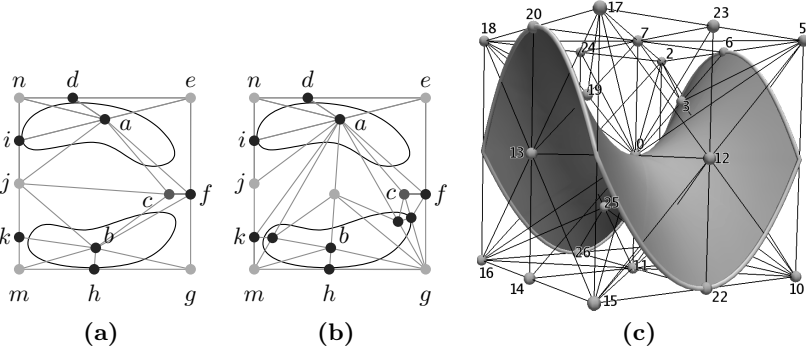
However, there is no guarantee that (C1) and (C2) are satisfied on all edges and faces of  $T_1$  that are not included (*stricto* or *lato sensu*) in  $M$ .

*Example 2.* Consider the scalar field defined by  $\Phi_2(x, y, z) = x^2 - y^2 + z$  over a single Lagrange Q2 element with linear geometry and coordinates in  $[-1, 1]^3$ , as illustrated in Fig. 3(c): each restriction of  $\Phi_2$  to the faces perpendicular to the  $z$ -axis has a critical point at the corresponding face center (labeled 7 and 11), and each restriction of  $\Phi_2$  to the edges perpendicular to the  $z$ -axis has a critical point at the corresponding edge midpoint (labeled 6, 23, 20, 24, 22, 14, and 26 for those that are visible), whereas  $\Phi_2$  proper does not have any critical point. Upon completion of TETRAHEDRALIZE-INTERIOR, all of these points have been inserted in  $T_1$  which contains, among others, edges from the element center to points 6, 20, 22, and 26. It is easy to check that the restrictions of  $\Phi_2$  to each of these edges have a critical point at the edge midpoint, and thus (C1) is not satisfied across  $T_1$ , that therefore requires further modification.

We must therefore examine how, in general,  $T_1$  can be modified so in order to satisfy (C1) and (C2). A natural question is to wonder whether it is possible to perform a series of edge-face flips on the  $T_1$  so that the final tessellation satisfies the desired properties.

*Example 3.* Given the triangulation of a face in Fig. 2(c) for which additional critical points (of the restrictions of the field to some of the new edges) have appeared, one can easily perform a series of edge flips so that the final connectivity satisfies (C1), as shown in the resulting tessellation in Fig. 3(a).

Nevertheless, it is unclear whether it is always possible to retrieve a tessellation that satisfies (C1) using only edge flips. Although this may be the case, we have not further explored this path, because the matter is more complicated as both (C1) and (C2) must be satisfied in a problem that is intrinsically three-dimensional: for instance, although any 2-D triangulation can always be converted to a Delaunay triangulation by a finite sequence of edge flips, but



**Fig. 3.** (a) A triangulation satisfying (C1) with edge flips from Fig. 2(c). (b) Another triangulation satisfying (C1) using CORRECT-TRIANGLE-TOPOLOGY and the same initial tessellation. (c) Topology-based tetrahedralization of a single Lagrange Q2 element with linear geometry for the field  $\Phi_2(x, y, z) = x^2 - y^2 + z$ .

this result does **not** extend to 3-D tetrahedralizations [2], making us skeptical of the flipping approach. We therefore took a different route, guided by Proposition 4, as not only  $T_1$  satisfies (C3), but we also know that will not be altered by subsequent subdivisions of  $T_1$ . Rather than attempting to identify a set of “problem” entities and prove that they may always be flipped into a satisfactory configuration, we introduce each critical point of a restriction of  $\Phi$  to a new entity into the tessellation. Any edges and faces created by this operation must then be examined as well for critical points of  $\Phi$  restricted to their respective domains. However, it is only necessary to focus on critical points of the restrictions of  $\Phi$  to the previously marked edges and faces, as others already satisfy (C1) and (C2), thanks to TRIANGULATE-BOUNDARIES.

CORRECT-TETRAHEDRAL-TOPOLOGY( $M, \Phi, S, T_1$ )

```

1  while  $S$  not empty
2      do Pop  $t$  from  $S$ 
3           $C \leftarrow \emptyset$ 
4          for  $e \leftarrow$  marked edges of  $t$ 
5              do Insert critical points of  $e$  into  $C$ 
6          for  $f \leftarrow$  marked faces of  $t$ 
7              do Insert critical points of  $f$  into  $C$ 
8          for  $c \leftarrow C$ 
9              do Find  $t \in T_1$  s. t.  $c \in t$ 
10             Remove  $t$  from  $T_1$  and  $S$ 
11              $U \leftarrow \text{STAR}^3(c, t)$ 
12             for  $t' \leftarrow U$ 
13                 do if MARK-TETRAHEDRON( $t'$ )
14                     then Push  $t'$  onto  $S$ 
15             Insert  $t'$  into  $T_1$ 

```

*Remark 5.* The same approach as that explained in Remark 4 is used here, using a tolerance to avoid refining below a certain size. Therefore, and for the same reason, this allows the algorithm to possibly terminate faster than normally possible if all critical points created on new edges and faces and the refinement proceeds had to be inserted without regard to the distance to an already inserted vertex.

**Proposition 5.** *If, for all  $e$  in  $M$ , all critical points of the restriction of  $\Phi_e$  to any arbitrary face are isolated, then CORRECT-TETRAHEDRAL-TOPOLOGY terminates. In addition, upon termination, (C1), (C2) and (C3) are satisfied.*

*Proof.* To establish this result, it is sufficient to make sure the algorithm terminates, starting from any arbitrary face of an arbitrary element in  $M$ . So, let  $f_i$  be one of the faces of an arbitrary  $e \in M$ , and we then shall prove that CORRECT-TETRAHEDRAL-TOPOLOGY terminates.

First, remark that if the restriction  $\Phi_{e|f_{ij}}$  of  $\Phi_e$  to one edge  $f_{ij}$  of  $f_i$  has a non-isolated critical point, then the derivative of  $\Phi_{e|f_{ij}}$  vanishes along a nonempty open segment of  $f_{ij}$ , and therefore has an infinity of zeros. Because this derivative is itself a univariate polynomial function, it can thus only be zero everywhere, and thus  $\Phi_{e|f_{ij}}$  is constant along the edge. Therefore, the only case when non-isolated critical points along a bounding edge of  $f_i$  arises is when the interpolant is constant along that edge, and therefore no other points than its endpoints are contained in  $P$ .  $P$  is indeed a finite set, as polynomials can only have a finite number of isolated critical points.

Now assume the restriction  $\Phi_{e|f_i}$  of  $\Phi_e$  to the interior of  $f_i$  has  $n \in \mathbb{N}^*$  critical points. The innermost loop of CORRECT-TETRAHEDRAL-TOPOLOGY will insert these  $n$  points, and yield a triangulation of  $f_i$  in  $N \in \mathbb{N}^*$  triangles  $t_{i,k}$ , such that  $\cup_{k=1}^N t_{i,k} = f_i$  and

$$(\forall 1 \leq k, k' \leq N) \quad k \neq k' \iff \overset{\circ}{t}_{i,k} \cap \overset{\circ}{t}_{i,k'} = \emptyset,$$

where all the points of  $C$  are vertices of some of  $t_{i,k}$  ( $\overset{\circ}{t}$  is the interior of  $t$  in the sense of the natural topology induced on  $t$  by embedding it in  $\mathbb{R}^2$ ). Therefore, none of the restrictions of  $\Phi_e$  to  $t_{i,k}$  has an internal critical point (otherwise this point would belong to  $C$ , which is impossible because all points of  $C$  are vertices of some of  $t_{i,k}$ ).

However, the restriction of  $\Phi_e$  to some edges of this triangulation of  $f_i$  may have critical points<sup>4</sup>. Denote  $\eta$  such an edge. If the restriction of  $\Phi$  to  $\eta$  has any non-isolated critical point, then the same argument as above holds and thus the corresponding edges do not need to be further refined. On the contrary, if such an edge critical point is isolated (in this case, the edge must be internal to  $f_i$ , as all isolated critical points along the edges of  $f_i$  have been inserted priorly), then CORRECT-TETRAHEDRAL-TOPOLOGY recursively proceeds on

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<sup>4</sup>In other words, the subdivision of  $f_i$  cannot create new face critical points, but it can create new edge critical points.

$\eta$ . However, the process terminates because all face critical points are supposed to be isolated according to the hypothesis. Therefore, for each critical point  $p_i$  of the restriction of  $\Phi_e$  to  $f_i$ , there exists a neighborhood of  $p_i$  in which all directional derivatives of  $\Phi_e$  are nonzero and thus, there exists a finite number of triangle subdivisions after which no edge critical points are left (because such a critical point implies one directional derivative is equal to 0).

Finally, upon completion of the algorithm, for the same reasons as for TETRAHEDRALIZE-INTERIOR, (C3) is satisfied for each tetrahedron of the final partition (the final tetrahedra are subdivisions of initial tetrahedra that all satisfied (C3), according to Proposition 4). $\square$

*Example 4.* Consider the same case as Example 2: the execution of CORRECT-TETRAHEDRAL-TOPOLOGY results in the insertion of the 4 previously mentioned edge midpoints (3 of which are visible in Fig. 3(c), labeled 3, 19, and 25) in a sub-tetrahedralization of  $T_1$  that becomes the final tessellation, across which conditions (C1) through (C4) are satisfied. In this case, only one level of refinement was necessary, as none of the edges and faces created upon insertion of the 4 aforementioned edge midpoints violates (C1) or (C2).

*Remark 6.* Note that the hypothesis of Proposition 5 is very stringent, as it is not limited to faces of the mesh, but extends to all possible faces. In fact, it is sufficient that the restrictions of  $\Phi_e$  to any face of the successive partitions of  $e$  only has isolated critical points. However, as this condition depends on the particular subdivision path, it is, albeit weaker, more difficult to prove.

### 3 Application to Isocontouring

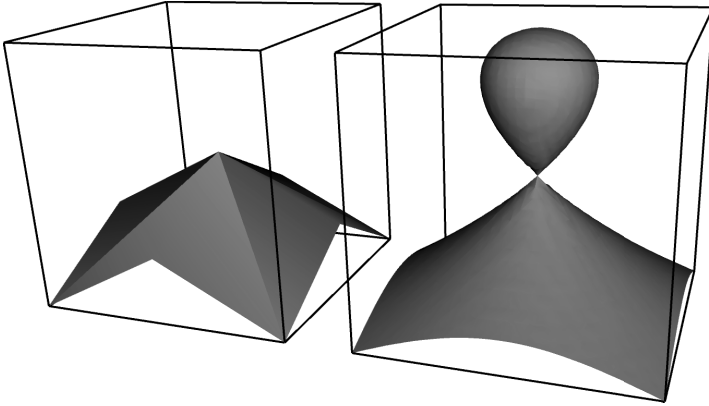
We now illustrate the application of our method to isocontouring, and the discuss an additional benefit of  $\kappa$ -tessellation that comes as a side-effect. Consider the following scalar field:

$$\begin{aligned} \Phi_3 : [-1, 1]^3 &\longrightarrow \mathbb{R} \\ (x, y, z)^T &\mapsto x^2 + y^2 + z^2(z - 1), \end{aligned}$$

interpolated over a single Q3 Lagrange element with linear geometry, i.e., a tricubic (hence with total degree 9) cell obtained by tensorization of cubic Lagrange interpolants. The test consists of representing the 0-isocontour of  $\Phi_3$ , which is tricky because an entire lobe of the resulting isosurface is contained within the element.

Figure 4, left, shows that if a linear isocontouring marching cubes technique is applied (after uniform subdivision of the hexahedron into 48 tetrahedra), then a substantial part of  $\Phi^{-1}(0)$  is missing. This example is interesting because the missing part of  $\Phi^{-1}(0)$  is not a disconnected component and, therefore, (C3) is *not* violated for this particular isovalue. While intuition may suggest that a linear isocontouring technique should retrieve the correct





**Fig. 4.** Tricubic isocontouring of  $\Phi_3$  for the isovalue 0, using a linear isocontouring technique (left), and using our approach (right).

topology of  $\Phi^{-1}(0)$  in the interior of the hexahedral element, in fact  $\Phi^{-1}(0)$  is not a 2-dimensional submanifold of  $\mathbb{R}^3$  at (and only at) point  $(0, 0, 0)$  (where  $\Phi^{-1}(0)$  is not simply connected). Therefore,  $\Phi^{-1}(0)$  is not a surface and this causes the isocontouring algorithm to fail, as illustrated. Moreover, the Implicit Function Theorem shows that for any value of  $\alpha$  in  $] -2, 0[$ ,  $\Phi^{-1}(\alpha)$  is a surface, but can also easily check that for such values of  $\alpha$ , it is not connected; in fact, it has 2 disjoint connected components, one of which is entirely contained in the interior of the element. Hence, in this case, (C3) is indeed violated; which also causes the linear isocontouring technique to fail (by missing this connected component).

From what we already know of our topology-based technique, we can expect it to properly retrieve  $\Phi^{-1}(\alpha)$  for these problematic values of  $\alpha$  in  $] -2, 0[$ , and it does indeed. However, what was not initially expected, is that it could fix the case when  $\alpha = 0$ , as we do not expect the technique to handle non-manifold isocontours. Nonetheless, as shown in Figure 4, right, our topology-based method is able to construct the correct topology of the non-simply connected  $\Phi^{-1}(0)$ . And in fact, this is not an anecdotal effect valid for this example only, but we can see that this is always the case: by “eliminating” critical points from the final tessellation, the method ensures that (thanks to the Implicit Function Theorem), the isocontour is locally a 2-dimensional manifold within the interior of each element. In other words, the scheme produces a tessellation that not only satisfies criteria (C1) through (C4), but additionally ensures that the isocontour is indeed a surface inside each element. Note that this extends to the lower-dimensional case, for the same reason: the isocontours inside the faces of the final tessellation are simply connected curves.

## 4 Conclusions

We have outlined an algorithm for partitioning finite elements to which is associated a scalar field into a  $\kappa$ -compatible tessellation, and proved that it works and terminates under a limited set of assumptions. This technique allows to easily adapt visualization and other post-processing tools to higher order elements. We have illustrated this methodology with the isocontouring operation.

Future work will include estimating the computational complexity of the algorithm as a function of input parameters such as the order of the interpolant, and improving multivariate polynomial system solution strategies.

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