

A Convergent Finite Volume type O-method on Evolving Surfaces

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Abstract. We present a finite volume scheme for anisotropic diffusion on evolving hypersurfaces. The underlying motion is assumed to be described by a fixed, not necessarily normal, velocity field. The ingredients of the numerical method are an approximation of the family of surfaces by a family of interpolating polygonal meshes, where grid vertices move on motion trajectories, a consistent finite volume discretization of the induced transport on the cells (polygonal patches), and a proper incorporation of a diffusive flux balance at polygonal faces. The main stability results and convergence estimate are obtained.

Keywords: Finite volume method, evolving surfaces, transport diffusion equations

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INTRODUCTION

In many applications, evolution problems do not reside on a flat Euclidean domain but on a curved hypersurface. Frequently this surface is itself evolving in time driven by some velocity field. In [1] Dziuk and Elliot proposed a finite element scheme for the numerical simulation of diffusion processes on such evolving surfaces, and in [2] a finite volume variant was proposed for simulation on simplicial meshes. In this paper we introduce a finite volume methodology for simulation of diffusion process on general evolving polygonal meshes. our finite volume is closely related to the pioneer work by C. Le Potier in [3] and the work of K. Lipnikov, M. Shashkov and I. Yotov in [4].

MATHEMATICAL MODEL

We consider a family of compact, smooth, and oriented hypersurfaces $\Gamma(t) \subset \mathbb{R}^n$ ($n = 2, 3$) for $t \in [0, t_{max}]$ generated by a time dependent function $\Phi : [0, t_{max}] \times \Gamma_0 \rightarrow \mathbb{R}^n$ defined on a reference surface Γ_0 with $\Phi(t, \Gamma_0) = \Gamma(t)$. Let us assume that Γ_0 is C^3 smooth and that $\Phi \in C^1([0, t_{max}], C^3(\Gamma_0))$. For simplicity we assume the reference surface Γ_0 to coincide with $\Gamma(0)$ (cf. 2).

We denote by $v = \partial_t \Phi$ the velocity of material points. The evolution of a conservative material quantity u with $u(t, \cdot) : \Gamma(t) \rightarrow \mathbb{R}$, which is propagated with the surface and simultaneously undergoes a linear diffusion on the surface, is governed by the parabolic equation

$$\dot{u} + u \nabla_\Gamma \cdot v - \nabla_\Gamma \cdot (\mathcal{D} \nabla_\Gamma u) = g \quad \text{on } \Gamma = \Gamma(t), \quad (1)$$

where $\dot{u} = \frac{d}{dt} u(t, x(t))$ is the (advective) material derivative of u , $\nabla_\Gamma \cdot v$ the surface divergence of the vector field v , $\nabla_\Gamma u$ the surface gradient of the scalar field u , g a source term with $g(t, \cdot) : \Gamma(t) \rightarrow \mathbb{R}$ and $\mathcal{D}$ a diffusion tensor on the tangent bundle. Here we assume a symmetric, uniformly coercive C^2 diffusion tensor field on whole $\mathbb{R}^n$ to be given, whose restriction on the tangent plane is then effectively incorporated in the model. With a slight misuse of notation, we denote this global tensor field also by $\mathcal{D}$. Furthermore, we impose an initial condition $u(0, \cdot) = u_0$ at time 0, and treat the case of surfaces with boundary. We then consider a Dirichlet boundary condition $u(t, \cdot) := u|_{\partial\Gamma(t)}$ on the boundary $\partial\Gamma(t)$. Let us assume that the mappings $(t, x) \rightarrow u(t, \Phi(t, x))$, $v(t, \Phi(t, x))$, and $g(t, \Phi(t, x))$ are $C^1([0, t_{max}], C^3(\Gamma_0))$, $C^0([0, t_{max}], C^3(\Gamma_0))$ and $C^1([0, t_{max}], C^1(\Gamma_0))$ regular, respectively.

DERIVATION OF THE FINITE VOLUME SCHEME

For the ease of presentation we restrict ourselves to the case of two dimensional surfaces in $\mathbb{R}^3$. let us give some preliminary definitions

Definition 0.1 (*Cell, cell center and vertices*) We call cell S a continuous 3D fan of triangles, where each triangle shares an edge with the preceding triangle, and all triangles share a common pivot point X_S called cell center or center point. The corner points p_i , ($i = 0, 1, \dots$) as depicted on Figure 1 (left) will be called vertices.

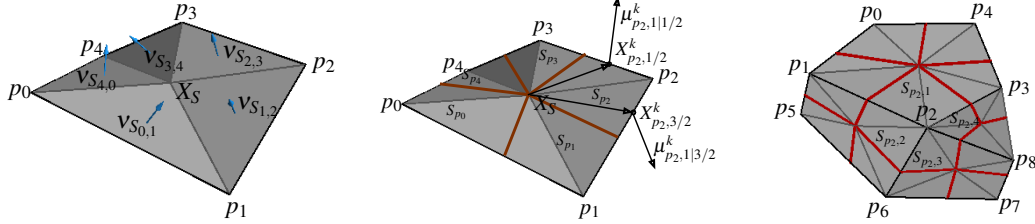


Figure 1. Admissible cell S and corresponding normals $v_{S_{i,j}}$ to triangles (left), sub-cells S_{p_i} of S , virtual unknowns $X_{p2,1/2}^k$, $X_{p2,3/2}^k$ of $S_{p2,1}$ and associated contravariant vectors $\mu_{p2,1/2}^k$, $\mu_{p2,3/2}^k$ (middle), sub-cells $S_{p2,j}$ around the vertex p_2 (right).

Definition 0.2 (*Admissible cell*) Let S be a cell, X_S its center point, and p_i ($i = 0, 1, \dots, n_S - 1$) its n_S vertices. For a given vertex p_i we denote by $v_{S_{i,j}} = \overrightarrow{X_S p_i} \wedge \overrightarrow{X_S p_j} / (\|\overrightarrow{X_S p_i} \wedge \overrightarrow{X_S p_j}\|)$ ($j \equiv (i+1) \bmod n_S$) the normal of the triangle $[X_S, p_i, p_j]$ if it has a non zero measure. We will then call S admissible cell if for any $i, j \equiv (i+1) \bmod n_S, l \equiv (j+1) \bmod n_S, m \in \{0, 1, \dots, n_S - 1\}$ $\|\overrightarrow{X_S p_i}\| \leq \max_{l,m} \|\overrightarrow{p_l p_m}\|$ and $v_{S_{i,j}} \cdot v_{S_{j,l}} > 0$ for well defined normals.

Definition 0.3 (*admissible polygonal surface*) We define an admissible polygonal surface Γ_h as a C^0 union of admissible cells where for two different cells $S_i, S_k \subset \Gamma_h$, $S_i \cap S_k$ is either an interface $\sigma = S_i|S_k = [p_i p_j]$, a vertex p_i , or is merely empty.

We now consider a sequence of admissible polygonal surfaces $\{\Gamma_h^k\}_{k=0, \dots, k_{max}}$ with Γ_h^k interpolating $\Gamma(t_k)$ for $t_k = k\tau$ and $k_{max}\tau = t_{max}$. Here, h indicates the maximal diameter of a cell on the whole sequence of polygonization, τ the time step size and k the index of a time step. All polygonisation share the same grid topology, and given the set of vertices p_j^0

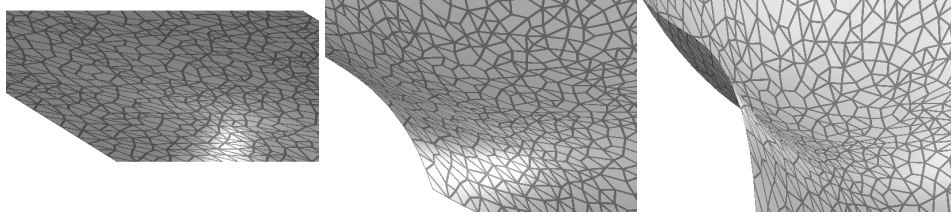


Figure 2. Zoom of sequence of polygonization Γ_h^k interpolating a surface in its evolution.

on the initial polygonal surface Γ_h^0 , the vertices of Γ_h^k lie on motion trajectories. Thus, they are evaluated based on the flux function Φ , i. e. $p_j(t_k) = \Phi(t_k, p_j^0)$. Upper indices denote the explicit geometric realization at the corresponding time step. We further assume that at each time step t_k , the Euclidean distance from any center point X_S^k to $\Gamma(t_k)$ is less than h^2 and the sub-cells $S_{i,j}^k := [X_S^k, p_i^k, p_{i+1}^k]$ ($i+1$ being the following index) are uniformly regular triangles. At each time step t_k , we consider a virtual subdivision of each cell S^k with n_S vertices into n_S polygonal sub-cells $\{S_{p_i}\}_{i=1, \dots, n_S}$ sharing X_S as depicted on Figure 1 (middle). This induces a subdivision of each edge $\sigma = [p_i, p_j] \subset \partial S$ into two sub-edges. Let us consider a vertex p_i and reorganize its surrounding cells S_j counter clockwise (cf. Figure 1 (right)). We call $\sigma_{p_i, j-1/2}$ and $\sigma_{p_i, j+1/2}$ the two sub-edges of S_j incident at p_i , and on each of these sub-edges, we respectively put the virtual unknowns $X_{p_i, j-1/2}^k$ and $X_{p_i, j+1/2}^k$. We recall that the index “ $j-1/2$ ” refers to the next sub-edge.

On $S_{p_i, j}$, sub-cell of S_j containing p_i , we define the covariant vectors $e_{p_i, j|j-1}^k := X_{p_i, j-1/2}^k - X_{S_j}^k$ and $e_{p_i, j|j+1}^k := X_{p_i, j+1/2}^k - X_{S_j}^k$ and their contravariant counter part $\mu_{p_i, j|j-1}^k$ and $\mu_{p_i, j|j+1}^k$ in $T_{p_i}^k := \text{Span}\{e_{p_i, j|j-1}^k, e_{p_i, j|j+1}^k\}$ such that

$\mu_{p_i,j|j-1}^k \cdot e_{p_i,j|j-1}^k = 1$, $\mu_{p_i,j|j-1}^k \cdot e_{p_i,j|j+1}^k = 0$, $\mu_{p_i,j|j+1}^k \cdot e_{p_i,j|j-1}^k = 0$ and $\mu_{p_i,j|j+1}^k \cdot e_{p_i,j|j+1}^k = 1$ (cf. Figure 1 (middle)). Using this dual system of vectors, we define on $S_{p_i,j}$ the natural approximation of tangential gradient $\nabla u(t_k, \cdot)|_{S_{p_i,j}} \approx \nabla_{p_i,j}^k u := [u(\mathcal{P}^k(X_{p_i,j-1/2}^k)) - u(\mathcal{P}^k(X_{p_i,j}^k))] \mu_{p_i,j|j-1}^k + [u(\mathcal{P}^k(X_{p_i,j+1/2}^k)) - u(\mathcal{P}^k(X_{p_i,j}^k))] \mu_{p_i,j|j+1}^k$. Here, $\mathcal{P}^k$ is the orthogonal projection onto the smooth surface when applied to interior points of Γ^k , and the orthogonal projection onto the boundary of the smooth surface when applied to boundary points of Γ^k . Based on these notational preliminaries, we can now derive a suitable finite volume discretization. Let us integrate (1) in $\{(t, x) | t \in [t_k, t_{k+1}], x \in S^{l,k}(t)\}$ ($S^{l,k}(t) := \Phi(t, \mathcal{P}^k(S^k)) \cap \Gamma(t_k)$) where $\mathcal{P}^k$ is taken to be the projection of onto a suitable extension of $\Gamma(t_k)$. We should mention here that $\mathcal{P}^k$ is context depend.

$$\int_{t_k}^{t_{k+1}} \int_{S^{l,k}(t)} g \, da \, dt \approx \tau m_S^{k+1} G_S^{k+1} \quad (2)$$

where $G_S^{k+1} = g(t_{k+1}, \mathcal{P}^{k+1} X_S^{k+1})$ and m_S^{k+1} the measure of S^{k+1} . The use of the Leibniz formula leads to the following approximation of the material derivative

$$\int_{t_k}^{t_{k+1}} \int_{S^{l,k}(t)} (\dot{u} + u \nabla_{\Gamma} \cdot v) \, da \, dt = \int_{S^{l,k}(t_{k+1})} u \, da - \int_{S^{l,k}(t_k)} u \, da \approx m_S^{k+1} u(t_{k+1}, \mathcal{P}^{k+1} X_S^{k+1}) - m_S^k u(t_k, \mathcal{P}^k X_S^k). \quad (3)$$

Furthermore, integrating the elliptic term again over the temporal evolution of a lifted cell and applying Gauss' theorem, we derive the following approximation:

$$\begin{aligned} \int_{t_k}^{t_{k+1}} \int_{S^{l,k}(t)} \nabla_{\Gamma} \cdot (\mathcal{D} \nabla_{\Gamma} u) \, da \, dt &= \int_{t_k}^{t_{k+1}} \int_{\partial S^{l,k}(t)} \mathcal{D} \nabla_{\Gamma} u \cdot \mu_{\partial S^{l,k}(t)} \, dl \, dt \\ &\approx \tau \sum_{p_i} \left(m_{p_i,j(S)+1/2}^{k+1} (\mathcal{D}_S^{k+1} \nabla_{p_i,j(S)}^{k+1} u) \cdot n_{p_i,j(S)|j(S)+1}^{k+1} + m_{p_i,j(S)-1/2}^{k+1} (\mathcal{D}_S^{k+1} \nabla_{p_i,j(S)}^{k+1} u) \cdot n_{p_i,j(S)|j(S)-1}^{k+1} \right) \end{aligned} \quad (4)$$

where $\mu_{\partial S^{l,k}(t)}$ is the unit outward co-normal on $\partial S^{l,k}(t)$ tangential to $\Gamma(t)$ and $m_{p_i,j(S)+1/2}^{k+1}$ the measure of $\sigma_{p_i,j(S)+1/2}^{k+1}$. The summation is done over the vertices p_i of the topological cell S , $j(S)$ denotes the reordered index number of S as a cell of the cluster of cells around p_i ; $\mathcal{D}_S^{k+1} := \mathcal{D}(t_{k+1}, X_S^{k+1})$, and $n_{p_i,j(S)|j(S)+1}^{k+1}$ respectively $n_{p_i,j(S)|j(S)-1}^{k+1}$ being defined as the unit outward co-normal vectors of $S_{p_i,j(S)}^k$. These last vectors belong to the plane $T_{p_i,j}^{k+1}$ and are respectively normal to $\sigma_{p_i,j(S)+1/2}^{k+1}$ and $\sigma_{p_i,j(S)-1/2}^{k+1}$. To close our equation, we impose a flux continuity on sub-edges which is numerically expressed as

$$\begin{aligned} &m_{p_i,j+1/2}^{k+1} \left[(U_{p_i,j+1/2}^{k+1} - U_{p_i,j}^{k+1}) \lambda_{p_i,j|j+1}^{k+1} + (U_{p_i,j-1/2}^{k+1} - U_{p_i,j}^{k+1}) \lambda_{p_i,j-1|j+1}^{k+1} \right] \\ &+ m_{p_i,j+1/2}^{k+1} \left[(U_{p_i,j+3/2}^{k+1} - U_{p_i,j+1}^{k+1}) \lambda_{p_i,j+2|j+1}^{k+1} + (U_{p_i,j+1/2}^{k+1} - U_{p_i,j+1}^{k+1}) \lambda_{p_i,j+1|j}^{k+1} \right] = 0. \end{aligned} \quad (5)$$

where $\lambda_{p_i,j|j+1}^{k+1} := \mathcal{D}_S^{k+1} \mu_{p_i,j|j+1}^{k+1} \cdot n_{p_i,j|j+1}^{k+1}$, $\lambda_{p_i,j-1|j+1}^{k+1} := \mathcal{D}_S^{k+1} \mu_{p_i,j|j-1}^{k+1} \cdot n_{p_i,j|j-1}^{k+1}$, and $\lambda_{p_i,j+1|j-1}^{k+1}$, $\lambda_{p_i,j|j-1}^{k+1}$ similarly defined. We denote by $U_{p_i,j}^{k+1}$ the cell center unknowns and $U_{p_i,j+1/2}^{k+1}$ the sub-edge unknowns. (5) gives a local relations $M_{p_i}^{k+1} \bar{U}_{p_i,j+1/2}^{k+1} = N_{p_i}^{k+1} \bar{U}_{p_i,j}^{k+1}$ between the vector of cell centers unknowns $\bar{U}_{p_i,j}^{k+1}$, and the vector of sub-edge unknowns $\bar{U}_{p_i,j+1/2}^{k+1}$ around vertices p_i . The behavior of the system depends on the matrices $M_{p_i}^{k+1}$, $N_{p_i}^{k+1}$ and

$$Q_{p_i,j}^{k+1} := \begin{pmatrix} m_{p_i,j+1/2}^{k+1} \lambda_{p_i,j|j+1}^{k+1} & m_{p_i,j-1/2}^{k+1} \lambda_{p_i,j+1|j-1}^{k+1} \\ m_{p_i,j+1/2}^{k+1} \lambda_{p_i,j-1|j+1}^{k+1} & m_{p_i,j-1/2}^{k+1} \lambda_{p_i,j|j-1}^{k+1} \end{pmatrix} \quad (6)$$

We then assume from now on that the sub-edge points are chosen such that $Q_{p_i,j}^{k+1}$ is positive definite and $(M_{p_i}^{k+1})^{-1} N_{p_i}^{k+1}$ is positive. It is then clear that an admissible polygonization must allow this set-up. It is worth mentioning here That if all incident angles at vertices are less than 180 degrees, a good choice of center point X_S^k of each cell S^k combined with a suitable optimization procedure around the vertices will always guarantee the above mentioned condition. The scheme described in [2] is then a good example and $Q_{p_i,j}^{k+1}$ is a diagonal matrix in that case. We recall (2), (3), and (4) to obtain the following equation on each cell

$$\begin{aligned} &m_S^{k+1} U_S^{k+1} - m_S^k U_S^k - \tau \sum_{p_i} m_{p_i,j(S)+1/2}^{k+1} \left[(U_{p_i,j(S)+1/2}^{k+1} - U_{p_i,j(S)}^{k+1}) \lambda_{p_i,j(S)|j(S)+1}^{k+1} + (U_{p_i,j(S)-1/2}^{k+1} - U_{p_i,j(S)}^{k+1}) \lambda_{p_i,j(S)-1|j(S)+1}^{k+1} \right] \\ &+ m_{p_i,j(S)-1/2}^{k+1} \left[(U_{p_i,j(S)+1/2}^{k+1} - U_{p_i,j(S)}^{k+1}) \lambda_{p_i,j(S)+1|j(S)-1}^{k+1} + (U_{p_i,j(S)-1/2}^{k+1} - U_{p_i,j(S)}^{k+1}) \lambda_{p_i,j(S)|j(S)-1}^{k+1} \right] = \tau m_S^{k+1} G_S^{k+1} \end{aligned} \quad (7)$$

which together with (5), the initial condition and the boundary condition give the finite volume scheme. Let us define the following discrete H_0^1 semi-norm.

Definition 0.4 (Discrete energy semi-norm). For any $U^k \in \mathcal{V}_h^k$ (set of constant function on cells), we define

$$\|U^k\|_{1,\Gamma_h^k}^2 := \sum_{S^k} \sum_{p_i^k \in S^k} \left(U_{p_i,j(S)+1/2}^k - U_{p_i,j(S)}^k, U_{p_i,j(S)-1/2}^k - U_{p_i,j(S)}^k \right) Q_{sym,p_i,j(S)}^k \left(U_{p_i,j(S)+1/2}^k - U_{p_i,j(S)}^k, U_{p_i,j(S)-1/2}^k - U_{p_i,j(S)}^k \right)^\top \quad (8)$$

where the sub-edge values $U_{p_{i,j(S)+1/2}}^k$ are defined by (5), and the boundary values are taken to be uniformly zero. $Q_{\text{sym}, p_{i,j(S)}}^k = (Q_{p_{i,j(S)}}^k + (Q_{p_{i,j(S)}}^k)^\top) / 2$ denotes the symmetric part of $Q_{p_{i,j(S)}}^k$ defined in (6).

We are now able to establish the main properties of the scheme.

Proposition 0.5 *The problem (7) has a unique solution.*

A PRIORI ESTIMATES

For simplicity in the analysis, we assume X_S^k in the convex hull of the vertices of S^k , as well as the following:

$$|\Upsilon^k(t_{k+1}, X_S^k) - X_S^{k+1}| \leq Ch\tau, \quad |\Upsilon^k(t_{k+1}, X_{p_{i,j+1/2}}^k) - X_{p_{i,j+1/2}}^{k+1}| \leq Ch\tau, \quad |\Upsilon^k(t_{k+1}, y_{p_{i,j+1/2}}^k) - y_{p_{i,j+1/2}}^{k+1}| \leq Ch\tau \quad (9)$$

where $y_{p_{i,j+1/2}}^k$ is the point that defines the sub-edge $\sigma_{p_{i,j+1/2}}^k$ together with p_i , $\Upsilon^k(t_{k+1}, X_S^k)$, $\Upsilon^k(t_{k+1}, X_{p_{i,j+1/2}}^k)$ and $\Upsilon^k(t_{k+1}, y_{p_{i,j+1/2}}^k)$ denote the points with the same barycentric coordinate in the convex hull of the image of the vertices of S^k through Φ .

Theorem 0.6 (Discrete $\mathbb{L}^\infty(\mathbb{L}^2), \mathbb{L}^2(\mathbb{H}^1)$ energy estimate). *Let $\{U^k\}_{k=1, \dots, k_{\max}}$ be the discrete solution of (7) for a given discrete initial data $U^0 \in \mathcal{V}_h^0$ and the homogenous boundary condition, then there exists a constant C depending solely on $t_{\max}$ such that*

$$\max_{k=1, \dots, k_{\max}} \|U^k\|_{\mathbb{L}^2(\Gamma_h^k)}^2 + \sum_{k=1}^{k_{\max}} \tau \|U^k\|_{1, \Gamma_h^k}^2 \leq C \left(\|U^0\|_{\mathbb{L}^2(\Gamma_h^0)}^2 + \tau \sum_{k=1}^{k_{\max}} \|G^k\|_{\mathbb{L}^2(\Gamma_h^k)}^2 \right) \quad (10)$$

Theorem 0.7 Discrete $\mathbb{H}^1(\mathbb{L}^2), \mathbb{L}^\infty(\mathbb{H}^1)$ energy estimate). *We assume the sub-matrices $Q_{p_{i,j}}^k$ (cf. (6)) to be symmetric for any sub-cell $S_{p_{i,j}}^k$ around p_i^k . Let $\{U^k\}_{k=1, \dots, k_{\max}}$ be the discrete solution of (7) for given discrete initial data $U^0 \in \mathcal{V}_h^0$ and the homogenous boundary condition, then there exists a constant C depending solely on $t_{\max}$ such that*

$$\sum_{k=1}^{k_{\max}} \tau \|\partial_t^\tau U^k\|_{\mathbb{L}^2(\Gamma_h^k)}^2 + \max_{k=1, \dots, k_{\max}} \|U^k\|_{1, \Gamma_h^k}^2 \leq C \left(\|U^0\|_{\mathbb{L}^2(\Gamma_h^0)}^2 + \|U^0\|_{1, \Gamma_h^0}^2 + \tau \sum_{k=1}^{k_{\max}} \|G^k\|_{\mathbb{L}^2(\Gamma_h^k)}^2 \right) \quad (11)$$

where $\partial_t^\tau U^k = \frac{U^k - U^{k-1}}{\tau}$ is defined as a difference quotient in time.

CONVERGENCE

Theorem 0.8 (Error estimate). *We define the piecewise constant error functional on Γ_h^k for $k = 1, \dots, k_{\max}$*

$$E^k := \sum_{S^k} \left(u^{-l}(t_k, X^k) - U_S^k \right) \chi_{S^k}$$

measuring the pull back $u^{-l}(t_k, \cdot) = u(t_k, \mathcal{P}^k(\cdot))$ of the continuous solution $u(t_k, \cdot)$ of (1) at time t_k and the finite volume solution U^k of (7). χ_{S^k} denotes the characteristic function of the cell S^k . Furthermore, let us assume that $\|E^0\|_{\mathbb{L}^2(\Gamma_h^0)} \leq Ch$, then the error estimate

$$\max_{k=1, \dots, k_{\max}} \|E^k\|_{\mathbb{L}^2(\Gamma_h^k)}^2 + \tau \sum_{k=1}^{k_{\max}} \|E^k\|_{1, \Gamma_h^k}^2 \leq C(h + \tau)^2 \quad (12)$$

holds for a constant C depending on the regularity assumptions and the time $t_{\max}$.

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