INBETWEENING AUTO-ANIMATION VIA FOKKER-PLANCK DYNAMICS AND THRESHOLDING

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(Communicated by Hao-Min Zhou)

Abstract. We propose an equilibrium-driven deformation algorithm (EDDA) to simulate the inbetweening transformations starting from an initial image to an equilibrium image, which covers images varying from a greyscale type to a colorful type on planes or manifolds. The algorithm is based on the Fokker-Planck dynamics on manifold, which automatically incorporates the manifold structure suggested by dataset and satisfies positivity, unconditional stability, mass conservation law and exponentially convergence. The thresholding scheme is adapted for the sharp interface dynamics and is used to achieve the finite time convergence. Using EDDA, three challenging examples, (I) facial aging process, (II) coronavirus disease 2019 (COVID-19) pneumonia invading/fading process, and (III) continental evolution process are computed efficiently.

1. Introduction. Inbetweening auto-animation is to automatically generate animations (motions) given starting and end images. The classical method for auto-animation uses detailed kinematic equations for each object in the starting images, which is precise but time consuming due to case by case c.f. [1, 9].

Instead of analyzing the detailed kinematic equation for each object, we aim to propose an efficient and universal algorithm for inbetweening auto-animation based on the Fokker-Planck dynamics on manifold and thresholding. We call this algorithm equilibrium-driven deformation algorithm (EDDA).

EDDA regards the end image as an equilibrium state of a Fokker-Planck equation and the inbetweening motion is driven by an underlying potential force determined by the equilibrium. This viewpoint is especially useful when the detailed physical

2020 Mathematics Subject Classification. Primary: 68U10, 94A08; Secondary: 37A30, 65M08.

Key words and phrases. Morphing, shape evolution, ergodicity, upwind scheme, Voronoi tessellation.

J.-G. Liu was supported in part by the National Science Foundation (NSF) under award DMS-1812573. G. Jin was supported in part by the the National Science Foundation of Guangdong Province under award 2019A1515011487, the China Postdoctoral Science Foundation under award 2020M681269 and the Fundamental Research Funds for the Central Universities under award 20184200031610059.

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process is not clear or hard to describe. For instance, the inbetweening motion of aging process, tumor growth, pneumonia invading for coronavirus disease 2019 (COVID-19) or the formation of current continents/oceans starting from Pangaea.

We first consider a Fokker-Planck equation in a flat domain $\Omega \subset \mathbb{R}^\ell$ with a unique equilibrium $\pi$ and no-flux boundary condition in Section 2 and then we propose an efficient solver for this Fokker-Planck equation in Section 3. The numerical solver for this part is based on structured grids and finite volume method [3]. An unconditionally stable explicit time discretization is introduced, which automatically enjoys positivity, mass conservation law, exponentially convergence and also efficiency. For a Fokker-Planck equation on a closed manifold, we propose a similar efficient solver based on point clouds and the associated Voronoi tessellation in Section 4. The Voronoi tessellation automatically gives the manifold information and can be used to approximate surface gradient/divergence in the Fokker-Planck equation. Based on this, an analogue unconditionally stable explicit time discretization is introduced.

To realize the end image (the equilibrium) at a finite time and the shape dynamics of the inbetweening motion, we combine the explicit-time-discretization of the Fokker-Planck equation with the thresholding dynamics. When the equilibrium image has a sharp interface, the scheme adapting thresholding step converges faster than the pure Fokker-Planck iteration and the relative error reaches machine accuracy at a finite time. In Section 3.2, we prove the finite-time convergence to the equilibrium for the Fokker-Planck solver with the thresholding adjustment by using a $\ell^1$ characterization for the thresholding adjustment.

In Section 5, we apply EDDA proposed for either structured grids on $\Omega \subset \mathbb{R}^\ell$, or for point-clouds which suggests an underlining manifold to conduct three challenging and important examples: (I) facial aging process, (II) COVID-19 pneumonia invading/fading, and (III) continental evolution process. In Example (I), inbetweening facial aging process at each time is simulated and potentially reveals the detailed changes of different part of human face over time. In Example (II), the inbetweening evolution of COVID-19 pneumonia invading before treatment and the fading away after treatment are simulated, which shows a good agreement with computerized tomography (CT) scans and also reveals promising application in the studying of pathology for COVID-19. In Example (III), the Fokker-Planck dynamics and thresholding adjustments are combined together to simulate the continental drifting process, which may suggest a new explanation for the formation of the current five continents of the world. From those examples from quite different research fields, EDDA are shown to be a very efficient and universal method with enormous potential applications in other fields of science and industry.

2. Fokker-Planck equation and equilibrium. Suppose $\Omega \subset \mathbb{R}^\ell$ is a bounded domain. Assume the end image on $\Omega$ is described by an equilibrium density $\pi(\mathbf{x}) : \Omega \rightarrow \mathbb{R}$. The value of $\rho$ indicates the gray level of the image for a grayscale image. In the case of Red-Green-Blue (RGB) image, we use three separate densities to indicate the RGB levels of the image separately. Then with $\pi \propto e^{-\phi}$, the Fokker-Planck equation is given by

\begin{equation}
\partial_t \rho = \Delta \rho + \nabla \cdot (\rho \nabla \phi) = \nabla \cdot \left( \pi \nabla \left( \frac{\rho}{\pi} \right) \right)
\end{equation}

with initial data $\rho_0$ satisfying

\begin{equation}
\int_\Omega \rho_0 \, d\mathbf{x} = \int_\Omega \pi \, d\mathbf{x}.
\end{equation}
We consider the following natural no-flux boundary condition
\begin{equation}
\mathbf{n} \cdot \nabla \left( \frac{\rho}{\pi} \right) = 0 \quad \text{on } \partial \Omega.
\end{equation}
See Section 4 for a Fokker-Planck equation on a $d$ dimensional smooth closed sub-manifold of $\mathbb{R}^d$.

Now we state the ergodicity of the Fokker-Planck equation (2.1). Assume
\begin{equation}
\pi > 0, \quad \pi \in C^1(\overline{\Omega}).
\end{equation}
Let $L^2(\Omega; \frac{1}{\pi} \, dx)$ be the weighted $L^2$ space. Define the Fokker-Planck operator for (2.1) as $L^* : D(L^*) \subset L^2(\Omega; \frac{1}{\pi} \, dx) \to L^2(\Omega; \frac{1}{\pi} \, dx)$ with $D(L^*) := \{ u \in H^2(\Omega; \frac{1}{\pi} \, dx); \partial_n \frac{u}{\pi} = 0 \text{ on } \partial\Omega \}$
\begin{equation}
L^* u := -\nabla \cdot \left( \pi \nabla \frac{u}{\pi} \right).
\end{equation}
This can be regarded as the adjoint operator of the generator $L = -\frac{1}{\pi} \nabla \cdot (\pi \nabla)$ of the Fokker-Planck equation (2.1). One can see $L^*$ is a self-adjoint operator in $L^2(\Omega; \frac{1}{\pi} \, dx)$ with compact resolvent $(\lambda I + L^*)^{-1}$ for $\lambda$ large enough. Thus $L^*$ has only discrete spectrum without finite accumulation points. Furthermore, since $\pi > 0$, for $\rho \in D(L^*)$,
\begin{equation}
L^* \rho = 0, \implies \int \pi |\nabla \frac{\rho}{\pi}|^2 \, dx = 0, \implies \rho = c\pi.
\end{equation}
Therefore, we conclude 0 is the simple principal eigenvalue of $L^*$ with the corresponding eigenfunction $\pi$, which leads to the spectral gap of $L^*$ in $L^2(\Omega; \frac{1}{\pi} \, dx)$, i.e.
\begin{equation}
\langle L^* u, u \rangle_{\frac{1}{\pi}} \geq c\langle u, u \rangle_{\frac{1}{\pi}}, \quad \text{for } u \text{ s.t. } \langle u, \pi \rangle_{\frac{1}{\pi}} = 0.
\end{equation}
Thus due to $\int (\rho - \pi) \, dx = 0$, we have the following Poincare's inequality
\begin{equation}
\int |\nabla \left( \frac{\rho}{\pi} - 1 \right)|^2 \pi \, dx \geq c \int \left( \frac{\rho}{\pi} - 1 \right)^2 \pi \, dx.
\end{equation}
Therefore, multiplying (2.1) by $\frac{\rho}{\pi} - 1$, by (2.3) we have
\begin{equation}
\frac{1}{2} \frac{d}{dt} \int \left( \frac{\rho}{\pi} - 1 \right)^2 \pi \, dx = -\int \pi \mid \nabla \frac{\rho}{\pi} \mid^2 \, dx \leq -c \int \left( \frac{\rho}{\pi} - 1 \right)^2 \pi \, dx,
\end{equation}
which gives the ergodicity that
\begin{equation}
\| \rho(\cdot, t) - \pi \|_{L^2(\Omega; \frac{1}{\pi} \, dx)} \leq e^{-ct} \| \rho(\cdot, 0) - \pi \|_{L^2(\Omega; \frac{1}{\pi} \, dx)}.
\end{equation}

3. EDDA based on structured grids. We present the numerical method based on structured grids for a Fokker-Planck equation on 2D domain $\Omega := [a, b] \times [c, d]$. Let the grid size be $\Delta x = \frac{b-a}{N}$, $\Delta y = \frac{d-c}{M}$. Define the cells as
\begin{equation}
C_{ij} = ((i-1)\Delta x, i\Delta x) \times ((j-1)\Delta y, j\Delta y), \quad i = 1, \ldots, N, \quad j = 1, \ldots, M.
\end{equation}
Then the cell centers $(x_i, y_j)$ are
\begin{equation}
x_i = a + \left( i - \frac{1}{2} \right) \Delta x, \quad y_j = c + \left( j - \frac{1}{2} \right) \Delta y, \quad i = 1, \ldots, N, \quad j = 1, \ldots, M.
\end{equation}
We use $\rho_{i,j}$ to approximate the value of $\rho(x_i, y_j)$ and take $\pi_{i,j} = \pi(x_i, y_j) > 0$. Then the continuous-time finite volume scheme is

\[
\frac{\partial \rho_{i,j}}{\partial t} = \frac{1}{\Delta x^2} \left( \frac{\rho_{i+1,j} + \rho_{i+1,j} - \rho_{i,j} - \rho_{i-1,j}}{2} \right) - \frac{\rho_{i,j} + \rho_{i,j} - \rho_{i-1,j} - \rho_{i-1,j}}{2} \frac{\pi_{i,j} + \pi_{i-1,j}}{\pi_{i,j} + \pi_{i-1,j}} \frac{\rho_{i,j} + \rho_{i,j} - \rho_{i-1,j} - \rho_{i-1,j}}{2} \frac{\pi_{i,j} + \pi_{i-1,j}}{\pi_{i,j} + \pi_{i-1,j}} \right)
\]

for $i = 1, \cdots, N$, $j = 1, \cdots, M$ with the no-flux boundary condition (2.3). Here $\dot{\rho}_{i,j}$ refers to time derivative of $\rho_{i,j}$. We assume the equilibrium $\pi$ satisfies

\[
\pi_{0,j} = \pi_{1,j}, \pi_{N+1,j} = \pi_{N,j} \quad j = 1, \cdots, M,
\]

\[
\pi_{i,0} = \pi_{i,1}, \pi_{i,M+1} = \pi_{i,M} \quad i = 1, \cdots, N,
\]

then the no-flux boundary condition (2.3) is reduced to

\[
\rho_{0,j} = \rho_{1,j}, \rho_{N+1,j} = \rho_{N,j} \quad j = 1, \cdots, M,
\]

\[
\rho_{i,0} = \rho_{i,1}, \rho_{i,M+1} = \rho_{i,M} \quad i = 1, \cdots, N.
\]

Denote $\rho_{i,j}^k$ as the value of $\rho$ at $t^k = k\Delta t$ with time step $\Delta t$. When applying to high resolution images, the convergence to the equilibrium is very slow for the classical explicit time discretization for (3.3) because the spectral gap is very small and the Courant-Friedrichs-Lewy (CFL) constraint. Hence we introduce the following unconditionally stable explicit time discretization for (3.3)

\[
\frac{\rho_{i,j}^{k+1} - \rho_{i,j}^k}{\Delta t} = \frac{1}{\Delta x^2} \left( \frac{\rho_{i,j}^{k+1} - \rho_{i,j}^k}{2} \frac{\pi_{i,j} + \pi_{i,j+1}}{\pi_{i,j}} - \frac{\rho_{i,j}^{k+1} - \rho_{i,j}^k}{2} \frac{\pi_{i-1,j} + \pi_{i,j}}{\pi_{i,j}} \right)
\]

for $i = 1, \cdots, N$, $j = 1, \cdots, M$ with the no-flux boundary condition (3.5).

We now further simplify (3.6) as

\[
\left(1 + \frac{\Delta t}{\Delta x^2} \left( \frac{\pi_{i,j} + \pi_{i,j+1}}{2\pi_{i,j}} + \frac{\pi_{i-1,j} + \pi_{i,j}}{2\pi_{i,j}} \right) \right) \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} = \frac{\rho_{i,j}^k}{\pi_{i,j}} + \frac{\Delta t}{\Delta x^2} \left( \frac{\pi_{i,j} + \pi_{i,j+1}}{2\pi_{i,j}} + \frac{\pi_{i-1,j} + \pi_{i,j}}{2\pi_{i,j}} \right) \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}}
\]

Define

\[
\lambda_{i,j} := \frac{1}{\Delta x^2} \left( \frac{\pi_{i,j} + \pi_{i,j+1}}{2\pi_{i,j}} + \frac{\pi_{i-1,j} + \pi_{i,j}}{2\pi_{i,j}} \right) > 0,
\]

(3.8)
which is always positive due to the positivity of $\pi_{i,j}$. Then (3.7) can be rewritten as

\[(1 + \Delta t \lambda_{i,j}) \rho_{i,j}^{k+1} = \rho_{i,j}^k + \Delta t \left( \frac{\pi_{i,j} + \pi_{i+1,j} \rho_{i+1,j}^k}{\pi_{i+1,j}} + \frac{\pi_{i-1,j} + \pi_{i,j} \rho_{i,j}^k}{\pi_{i-1,j}} \right) \]

\[+ \Delta t \left( \frac{\pi_{i,j} + \pi_{i+1,j} \rho_{i+1,j}^k}{\pi_{i+1,j}} + \frac{\pi_{i,j-1} + \pi_{i,j} \rho_{i,j-1}^k}{\pi_{i,j-1}} \right). \]  

(3.9)

Denote $h := \max\{\Delta x, \Delta y\}$. From (3.9), we recast the scheme using a rescaled generator operator

\[
\frac{\rho_{i,j}^{k+1}}{\rho_{i,j}^k} - \frac{h^2}{1 + \Delta t \lambda_{i,j} \Delta x^2} \left( \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} - \frac{\rho_{i+1,j}^k}{\pi_{i+1,j}} \right) \]

\[= \frac{\Delta t}{2 \Delta x^2 \pi_{i,j}} \left[ \frac{\pi_{i,j} + \pi_{i+1,j}}{\pi_{i+1,j}} \left( \rho_{i+1,j}^k - \rho_{i,j}^k \right) - \frac{\pi_{i,j} + \pi_{i-1,j}}{\pi_{i-1,j}} \left( \rho_{i,j}^k - \rho_{i-1,j}^k \right) \right] \]

\[+ \frac{\pi_{i,j+1} + \pi_{i,j}}{2 \Delta y^2 \pi_{i,j}} \left( \rho_{i,j+1}^k - \rho_{i,j}^k \right) - \frac{\pi_{i,j} + \pi_{i,j-1}}{2 \Delta y^2 \pi_{i,j}} \left( \rho_{i,j}^k - \rho_{i,j-1}^k \right) \]

\[= -\Delta t (L_h \frac{\rho_{i,j}^k}{\pi_{i,j}}). \]

Now we state the the positivity, maximal principle, mass conservation law and ergodicity of the scheme (3.7) as follows.

**Proposition 1.** Let $\pi_{i,j} = \pi(x_i, y_j) > 0$. Let $\Delta t$ be the time step and consider the explicit scheme (3.6) for the numerical solution $\rho_{i,j}^k$ with boundary condition (3.5). Assume the initial data $\rho^0 > 0$ satisfies

\[
\sum_{i=1}^N \sum_{j=1}^M (1 + \Delta t \lambda_{i,j}) \rho_{i,j}^0 = \sum_{i=1}^N \sum_{j=1}^M (1 + \Delta t \lambda_{i,j}) \pi_{i,j}. \]  

(3.11)

Then we have

(i) positivity preserving property

\[
\rho_{i,j}^k > 0, \quad i = 1, \cdots, N, \quad j = 1, \cdots, M \quad \Rightarrow \quad \rho_{i,j}^{k+1} > 0, \]

(3.12)

\[i = 1, \cdots, N, \quad j = 1, \cdots, M; \]

(ii) the mass-conversation law

\[
\sum_{i=1}^N \sum_{j=1}^M (1 + \Delta t \lambda_{i,j}) \rho_{i,j}^{k+1} = \sum_{i=1}^N \sum_{j=1}^M (1 + \Delta t \lambda_{i,j}) \rho_{i,j}^k. \]

(3.13)

(iii) the unconditional maximal principle for $\frac{\rho_{i,j}^{k+1}}{\pi_{i,j}}$ in $\frac{\rho_{i,j}^k}{\pi_{i,j}}$

\[
\max_{i,j} \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} \leq \max_{i,j} \frac{\rho_{i,j}^k}{\pi_{i,j}}; \]

(3.14)

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(iv) the $l^\infty$ contraction

\[
\max_{i,j} \left| \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} - 1 \right| \leq \max_{i,j} \left| \frac{\rho_{i,j}^k}{\pi_{i,j}} - 1 \right| ;
\]

(v) the exponential convergence

\[
\left\| \left( \frac{\rho_{i,j}^k}{\pi_{i,j}} \right) - u^* \right\|_F \leq c\mu_2^k, \quad |\mu_2| < 1,
\]

where $\mu_2$ is the second eigenvalue of $A$ defined in (3.26), $u_{i,j}^* \equiv 1$ and $\| \cdot \|_F$ is the Frobenius norm of matrix.

Proof. For (i), from (3.9), since $\pi > 0$, we know $\rho_{i,j}^k > 0$ implies $\rho_{i,j}^{k+1} > 0$.

To prove (ii), taking summation in (3.9), we have the

\[
\pi > \sum_{i=1}^N \sum_{j=1}^M (1 + \Delta t \lambda_{i,j}) \rho_{i,j}^k
\]

\[
= \sum_{i=1}^N \sum_{j=1}^M \rho_{i,j}^k + \sum_{i=1}^N \sum_{j=1}^M \Delta t \frac{\pi_{i,j} + \pi_{i+1,j}}{2} \frac{\rho_{i+1,j}^k}{\pi_{i+1,j}} + \sum_{i=1}^N \sum_{j=1}^M \Delta t \frac{\pi_{i-1,j} + \pi_{i,j}}{2} \frac{\rho_{i-1,j}^k}{\pi_{i-1,j}}
\]

\[
+ \sum_{i=1}^N \sum_{j=1}^M \Delta t \frac{\pi_{i,j} + \pi_{i,j+1}}{2} \frac{\rho_{i,j+1}^k}{\pi_{i,j+1}} + \sum_{i=1}^N \sum_{j=1}^M \Delta t \frac{\pi_{N,j} + \pi_{N+1,j}}{2} \frac{\rho_{N+1,j}^k}{\pi_{N+1,j}}
\]

\[
- \sum_{j=1}^M \Delta t \frac{\pi_{0,j} + \pi_{1,j}}{2} \frac{\rho_{1,j}^k}{\pi_{1,j}} + \Delta t \frac{\rho_{N+1,j}^k - \rho_{1,j}^k}{\pi_{N+1,j}} ,
\]

where we used the no-flux boundary condition (3.4). Similarly, the third term in the RHS of (3.17) is

\[
\sum_{i=1}^N \sum_{j=1}^M \Delta t \frac{\pi_{i-1,j} + \pi_{i,j}}{2} \frac{\rho_{i-1,j}^k}{\pi_{i-1,j}} + \sum_{j=1}^M \Delta t \frac{\pi_{N,j} + \pi_{N+1,j}}{2} \frac{\rho_{N+1,j}^k}{\pi_{N+1,j}}
\]

\[
+ \sum_{j=1}^M \Delta t \frac{\pi_{0,j} + \pi_{1,j}}{2} \frac{\rho_{0,j}^k}{\pi_{0,j}}
\]
\[
= \sum_{i=1}^{N} \sum_{j=1}^{M} \frac{\Delta t}{\Delta x^2} \pi_{i+1,j} \pi_{i,j} \frac{\rho_{i,j}^k}{\pi_{i,j}} \frac{\Delta t}{\Delta x^2} \left( \rho_{N,j}^k - \rho_{0,j}^k \right).
\]

One can shift index for the last two terms in the RHS of (3.17) similarly. Therefore, using the no-flux boundary condition (3.5), we have the mass balance

\[
\sum_{i=1}^{N} \sum_{j=1}^{M} (1 + \Delta t \lambda_{i,j}) \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} = \sum_{i=1}^{N} \sum_{j=1}^{M} (1 + \Delta t \lambda_{i,j}) \frac{\rho_{i,j}^k}{\pi_{i,j}}.
\]

To prove (iii), directly taking maximum in the RHS of (3.7) implies

\[
(1 + \Delta t \lambda_{i,j}) \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} \leq (1 + \Delta t \lambda_{i,j}) \max_{i,j} \frac{\rho_{i,j}^k}{\pi_{i,j}},
\]

which leads to (4.16).

To prove (iv), subtract \((1 + \Delta t \lambda_{i,j})\) from both sides of (3.7) and then multiply by \(\text{sgn} \left( \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} - 1 \right)\). Thus using same argument with (iii), we have

\[
(1 + \Delta t \lambda_{i,j}) \left| \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} - 1 \right| \leq (1 + \Delta t \lambda_{i,j}) \max_{i,j} \left| \frac{\rho_{i,j}^k}{\pi_{i,j}} - 1 \right|,
\]

which implies (4.17).

Now we prove (v). Recall (3.10), i.e.

\[
\frac{\rho_{i,j}^{k+1}}{\pi_{i,j}} - \frac{\rho_{i,j}^k}{\pi_{i,j}} = -\Delta t \left( L_h \frac{\rho_{i,j}^k}{\pi} \right)_{i,j}
\]

\[
= \frac{\Delta t}{1 + \Delta t \lambda_{i,j}} \left[ \frac{\pi_{i,j}}{\Delta x^2 \pi_{i,j}} \left( \frac{\rho_{i+1,j}^k}{\pi_{i+1,j}} - \frac{\rho_{i,j}^k}{\pi_{i,j}} \right) - \frac{\pi_{i,j} + \pi_{i-1,j}}{\Delta x^2 \pi_{i,j}} \left( \frac{\rho_{i,j}^k}{\pi_{i,j}} - \frac{\rho_{i-1,j}^k}{\pi_{i-1,j}} \right) \right.
\]

\[
+ \frac{\pi_{i,j+1} + \pi_{i,j}}{\Delta y^2 \pi_{i,j}} \left( \frac{\rho_{i,j+1}^k}{\pi_{i,j+1}} - \frac{\rho_{i,j}^k}{\pi_{i,j}} \right) - \frac{\pi_{i,j} + \pi_{i,j-1}}{\Delta y^2 \pi_{i,j}} \left( \frac{\rho_{i,j}^k}{\pi_{i,j}} - \frac{\rho_{i,j-1}^k}{\pi_{i,j-1}} \right). \]

By shifting index and no-flux boundary condition (3.5) we have

\[
\langle -L_h \frac{\rho_{i,j}^k}{\pi}, \rho^k \rangle_{1+\Delta t \lambda} = -\Delta t \left[ \frac{\pi_{i,j} + \pi_{i+1,j}}{2 \Delta x^2 \pi_{i,j}} \left( \frac{\rho_{i+1,j}^k}{\pi_{i+1,j}} - \frac{\rho_{i,j}^k}{\pi_{i,j}} \right)^2 + \frac{\pi_{i,j} + \pi_{i-1,j}}{2 \Delta x^2 \pi_{i,j}} \left( \frac{\rho_{i,j}^k}{\pi_{i,j}} - \frac{\rho_{i-1,j}^k}{\pi_{i-1,j}} \right)^2 \right.
\]

\[
+ \frac{\pi_{i,j+1} + \pi_{i,j}}{2 \Delta y^2 \pi_{i,j}} \left( \frac{\rho_{i,j+1}^k}{\pi_{i,j+1}} - \frac{\rho_{i,j}^k}{\pi_{i,j}} \right)^2 + \frac{\pi_{i,j} + \pi_{i,j-1}}{2 \Delta y^2 \pi_{i,j}} \left( \frac{\rho_{i,j}^k}{\pi_{i,j}} - \frac{\rho_{i,j-1}^k}{\pi_{i,j-1}} \right)^2 \right],
\]

where the shorthand notation \(\langle \cdot , \cdot \rangle_{1+\Delta t \lambda}\) means the weighted inner product with the weight \(1 + \Delta t \lambda_{i,j}\). From this, we know

\[
\langle L_h \rho_{i,j}^\infty, \rho^\infty \rangle_{1+\Delta t \lambda} = 0 \quad \Rightarrow \quad \rho^\infty = c \pi = \pi,
\]

where \(c = 1\) due to the mass conservation law.
Denote \( u_{i,j}^k = \frac{\rho_{i,j}^k}{\pi_{i,j}} \). Then (3.10) is recast as
\[
\begin{align*}
  u_{i,j}^{k+1} &= \frac{1}{1 + \Delta t \lambda_{i,j}} \left[ u_{i,j}^k + \Delta t \left( \frac{\pi_{i,j} + \pi_{i+1,j} u_{i+1,j}^k}{2\Delta x^2 \pi_{i,j}} + \frac{\pi_{i,j} + \pi_{i-1,j} u_{i-1,j}^k}{2\Delta x^2 \pi_{i,j}} \right) 
  + \Delta t \left( \frac{\pi_{i,j+1} + \pi_{i,j} u_{i,j+1}^k}{2\Delta y^2 \pi_{i,j}} + \frac{\pi_{i,j} + \pi_{i,j-1} u_{i,j-1}^k}{2\Delta y^2 \pi_{i,j}} \right) \right] =: (Au^k)_{i,j}.
\end{align*}
\]

For presentation simplicity, we can regard \( u^k \) as a vector. By the Perron-Frobenius theorem, \( \mu_1 = 1 \) is the simple, principal eigenvalue of \( A \) with the ground state \( u_{i,j}^* \equiv 1 \) and other eigenvalues \( \mu_i \) of \( A \) satisfy \( |\mu_i| < \mu_1 \). Notice also the mass conservation for initial data \( u^0 = \rho^0 \) satisfying (4.15), i.e.,
\[
\langle u^0 - u^*, u^* \rangle (1 + \Delta t \lambda) = 0.
\]

Since also \( A \) is self-adjoint operator in the weighted space \( l^2((1 + \Delta \lambda)\pi) \), we can express \( u^0 \) using
\[
\begin{align*}
  u^0 - u^* &= \sum_{\ell=2}^{MN} c_{\ell} u_{\ell}, \quad u_{\ell} \text{ is the eigenfunction corresponding to } \mu_{\ell}.
\end{align*}
\]

Therefore, we have
\[
\begin{align*}
  u^k - u^* &= A^k(u^0 - u^*) = \sum_{\ell=2}^{MN} c_{\ell} \mu_{\ell}^k u_{\ell},
\end{align*}
\]
which concludes
\[
\|u^k - u^*\|_{F} \leq c|\mu_2|^k \quad \text{with } |\mu_2| < 1,
\]
where \( \| \cdot \|_{F} \) is the Frobenius norm. \( \square \)

**Remark 1.** In order to adjust the initial distribution \( \rho^0 \) such that the mass conservation law (3.11) is satisfied. We multiply \( \rho^0 \) by a constant \( c\text{adj} := \frac{\sum_{i,j}(1 + \Delta t \lambda_{i,j})\pi_{i,j}}{\sum_{i,j}(1 + \Delta t \lambda_{i,j})\rho_{i,j}^0} \), which does not affect the Fokker-Planck dynamics for \( \rho^k \). The image structure, as well as the inbetweening transformation, will remain the same after adjustment. To recover the original color scalar, one can use different interpolation models as a post-process, for instance the simple linear interpolation or using a Wasserstein–Fisher–Rao metric as for the unbalanced optimal transport.

**Lemma 3.1 (Weighted \( \ell^1 \)-contraction).** Let \( \rho_{i,j}^{k+1} \) be the solution obtained in (3.7). Then we have the weighted \( \ell^1 \)-contraction
\[
\begin{align*}
  \sum_{i,j} (1 + \Delta t \lambda_{i,j}) |\rho_{i,j}^{k+1} - \pi_{i,j}| \leq \sum_{i,j} (1 + \Delta t \lambda_{i,j}) |\rho_{i,j}^k - \pi_{i,j}|.
\end{align*}
\]

**Proof.** From (3.9), we know
\[
\begin{align*}
  (1 + \Delta t \lambda_{i,j}) (\rho_{i,j}^{k+1} - \pi_{i,j}) &= (\rho_{i,j}^k - \pi_{i,j}) + \frac{\Delta t}{\Delta x^2} \left( \frac{\pi_{i,j} + \pi_{i+1,j} \rho_{i+1,j}^k - \pi_{i+1,j}}{\pi_{i+1,j}} + \frac{\pi_{i,j} + \pi_{i-1,j} \rho_{i-1,j}^k - \pi_{i-1,j}}{\pi_{i-1,j}} \right) 
  + \frac{\Delta t}{\Delta y^2} \left( \frac{\pi_{i,j} + \pi_{i,j+1} \rho_{i,j+1}^k - \pi_{i,j+1}}{\pi_{i,j+1}} + \frac{\pi_{i,j} + \pi_{i,j-1} \rho_{i,j-1}^k - \pi_{i,j-1}}{\pi_{i,j-1}} \right).
\end{align*}
\]
Then taking the absolute value for both sides, we have

\[
\sum_{i,j} (1 + \Delta t \lambda_{i,j}) |\rho^{k+1}_{i,j} - \pi_{i,j}| \\
\leq \sum_{i,j} |\rho^k_{i,j} - \pi_{i,j}| \\
+ \sum_{i,j} \frac{\Delta t}{\Delta x^2} \left( \frac{\pi_{i,j} + \pi_{i+1,j}}{2} |\rho^{k}_{i+1,j} - \pi_{i+1,j}| + \frac{\pi_{i-1,j} + \pi_{i,j}}{2} |\rho^{k}_{i-1,j} - \pi_{i-1,j}| \right) \\
+ \sum_{i,j} \frac{\Delta t}{\Delta y^2} \left( \frac{\pi_{i,j} + \pi_{i,j+1}}{2} |\rho^{k}_{i,j+1} - \pi_{i,j+1}| + \frac{\pi_{i,j-1} + \pi_{i,j}}{2} |\rho^{k}_{i,j-1} - \pi_{i,j-1}| \right) \\
= \sum_{i,j} (1 + \Delta t \lambda_{i,j}) |\rho^k_{i,j} - \pi_{i,j}|,
\]

where the last equality comes from the same argument as the conservation law of total mass (3.20).

3.1. Thresholding adjustment for shape dynamics. In this section, we combine the thresholding scheme with the Fokker-Planck dynamics to generate the inbetweening motions with sharp interface, i.e., the density is described by linear combinations of two characteristic functions. In the computations for the continental evolution, one will see that the thresholding scheme also helps to achieve the finite time convergence to the sharp equilibrium density. The analytic result for the finite time convergence property will be proved in Section 3.2.

Notice the dynamics of the Fokker-Planck equation is invariant when replacing \( \rho \) by \( c \rho \). Therefore, the initial density shall be adjusted based on the mass conservation law (3.11); see Remark 1. After this initial adjustment, assume initial data \( \rho^0_{i,j} \in \{\rho^0_s, \rho^0_b\} \), which takes alternatively the value \( \rho^0_s, \rho^0_b \). Assume the equilibrium is \( \pi_{i,j} \in \{\pi_s, \pi_b\} \) which takes alternatively the value \( \pi_s, \pi_b \).

To combine the thresholding scheme with the Fokker-Planck dynamics, we need to choose the threshold constant \( \xi^k \) at each step to conserve (3.13) as follows:

**Step 1.** Given \( \rho^k_{i,j} \in \{\pi_s, \pi_b\} \), compute the explicit Fokker-Planck scheme (3.7) to update \( \tilde{\rho}^{k+1}_{i,j} \) for any \( i = 1, 2, \cdots, N \) and \( j = 1, 2, \cdots, M \).

**Step 2.** Choose thresholding constant \( \xi^{k+1} \) and define

\[
\rho^{k+1}_{i,j} := \pi_s \chi_{\{i,j; \rho^k_{i,j} \leq \xi^{k+1}\}} + \pi_b \chi_{\{i,j; \rho^k_{i,j} > \xi^{k+1}\}}
\]

such that \( \rho^{k+1} \) satisfies (3.13).

In Step 2, the constant \( \xi^{k+1} \) can be found using bisection such that

\[
f(\xi^{k+1}) := \sum_i (1 + \lambda_{i,j} \Delta t) \rho^{k+1}_{i,j} - \sum_i (1 + \lambda_{i,j} \Delta t) \pi_{i,j} = 0.
\]

3.2. Finite time convergence to equilibrium for the thresholding adjustment. In this section, we prove the finite time convergence to equilibrium for the thresholding adjustment combined with the classical explicit scheme. We first present the explicit scheme by changing the RHS of (3.6) to \( \rho^k \), i.e.,

\[
\frac{\rho^{k+1}_{i,j} - \rho^k_{i,j}}{\Delta t}
\]
where the constant we can always adjust $\tilde{\rho}_{i,j}$.

Without loss of generality, we assume

$$\sum_{i,j} \left(\frac{\rho_{i,j}^k}{\pi_{i,j}} - \frac{\rho_{i,j}^{k+1}}{\pi_{i,j}}\right)$$

for $i = 1, \ldots, N, j = 1, \ldots, M$ with the no-flux boundary condition (3.5). The CFL condition for this explicit scheme is $\max_{i,j} \lambda_{i,j} \Delta t < 1$. We have the mass conservation law

$$\sum_{i,j} \rho_{i,j}^k = \sum_{i,j} \rho_{i,j}^0$$

and the $\ell^1$-contraction under CFL condition

$$\sum_{i,j} |\rho_{i,j}^{k+1} - \pi_{i,j}| \leq \sum_{i,j} |\rho_{i,j}^k - \pi_{i,j}|.$$

For this explicit scheme (3.35), the associated thresholding adjustment for the sharp interface dynamics in Section 3.1 becomes

$$\rho_{i,j}^{k+1} := \pi_{i,j} + \xi \chi_{\{\pi_{i,j}^{k+1} \leq \xi \varepsilon^{k+1}\}} + \pi_{i,j} \chi_{\{\pi_{i,j}^{k+1} > \xi \varepsilon^{k+1}\}},$$

where the constant $\xi \varepsilon^{k+1}$ is chosen such that $\rho_{i,j}^{k+1} \in A := \{v_{i,j} \in \{\pi_{s}, \pi_{b}\}; \sum_{i,j} v_{i,j} = \sum_{i,j} \pi_{i,j}\}$.

Next, we claim the thresholding adjustment (3.38) is a minimizer in the subset $A$. To prove it, we first give an elementary rearrangement inequality.

**Lemma 3.2.** Let $a \leq b$, for any two numbers $u$ and $v$, we have the following elementary rearrangement inequality

$$|a - \min(u, v)| + |b - \max(u, v)| \leq |a - u| + |b - v|.$$

As a consequence, we have the following characterization for the thresholding adjustment

**Lemma 3.3.** For each $k$, let $\tilde{\rho}_{i,j}^k$ be the solution obtained from Fokker-Planck solver (3.35). Then the thresholding adjustment $\rho_{i,j}^k$ satisfies

$$\sum_{i,j} |\rho_{i,j}^k - \tilde{\rho}_{i,j}^k| = \min_{v \in A} \sum_{i,j} |v_{i,j} - \tilde{\rho}_{i,j}^k|.$$

**Proof.** Without loss of generality, we assume $\tilde{\rho}_{i,j}^k$ are distinct for any $i, j$. Otherwise, we can always adjust $\tilde{\rho}_{i,j}^k$ with a $\varepsilon$-perturbation such that the new $\tilde{\rho}_{i,j}^k$ is distinct and then take $\varepsilon \to 0$.

Thus there exists a constant $\xi^k$ such that $\rho_{i,j}^k \in A$ satisfies (3.38). Denote the points such that $\rho_{i,j}^k > \xi^k$ as $P^+ := \chi\{i, j; \rho_{i,j}^k > \xi^k\}$ whose cardinality is denoted as $n_0$. Then the cardinality of points such that $\rho_{i,j}^k \leq \xi^k$, denoted as $P^-$, is $n - n_0$.

Then we claim $\rho_{i,j}^k$ satisfies (3.40). Indeed, for any $v \neq \rho_{i,j}^k, v \in A$, it can be obtained from $\tilde{\rho}_{i,j}^k$ using (3.38) by switching $\ell$-pairs of points in $P^+$ and in $P^-$ for some integer $\ell \geq 1$. Then for such $\ell$-pairs of points, we apply Lemma 3.2 to conclude that switch of pairs in $P^+$ and $P^-$ leads to a larger $\ell^1$ norm

$$\sum_{i,j} |\rho_{i,j}^k - \tilde{\rho}_{i,j}^k| \leq \sum_{i,j} |v_{i,j} - \tilde{\rho}_{i,j}^k|.$$
Now we give the finite time convergence to equilibrium after applying the thresholding adjustment to the explicit Fokker-Planck scheme (3.35) provided the $\ell^1$ relative error is smaller than $\pi_b - \pi_s$. This proposition also explains the finite time convergence property in the continental evolution example in Section 5.3.

**Proposition 2.** Assume equilibrium $\pi_{i,j} \in \{\pi_s, \pi_b\}$. Let $\tilde{\rho}_{i,j}^k$ be the solution obtained from Fokker-Planck solver (3.35) and the thresholding adjustment $\rho_{i,j}^k$ be defined in (3.38). If the $\ell^1$ relative error reaches

$$\sum_{i,j} |\tilde{\rho}_{i,j}^k - \pi_{i,j}| < \pi_b - \pi_s$$

for some $k$ large enough, then the thresholding adjustment $\rho_{i,j}^k = \pi_{i,j}$.

**Proof.** Assume there exists $k$ such that (3.41). We prove $\rho_{i,j}^k = \pi_{i,j}$ by contradiction argument. If not, then

$$2(\pi_b - \pi_s) \leq \sum_{i,j} |\rho_{i,j}^k - \pi_{i,j}| \leq \sum_{i,j} |\rho_{i,j}^k - \tilde{\rho}_{i,j}^k| + \sum_{i,j} |\tilde{\rho}_{i,j}^k - \pi_{i,j}|$$

$$\leq 2 \sum_{i,j} |\tilde{\rho}_{i,j}^k - \pi_{i,j}| < 2(\pi_b - \pi_s),$$

where we used (3.40) in the third inequality. This is a contradiction and we conclude $\rho_{i,j}^k = \pi_{i,j}$. 

Based on the exponential convergence of $\tilde{\rho}$ to the equilibrium in Proposition (1), we know the condition (3.41) is satisfied after finite iteration steps. We also note that $\pi$ is the equilibrium for both the Fokker-Planck solver and the thresholding adjustment. Hence once $\rho$ reaches $\pi$, it will stay invariant. This proposition for the thresholding adjustment on structured grids provides analytic explanations for the finite time convergence of the continental evolution example in Section 5.3.

4. **EDDA based on point-clouds: Fokker-Planck equation on $N$.** Suppose $(N, d_N)$ is a $d$ dimensional smooth closed submanifold of $\mathbb{R}^\ell$. Assume the end image on $N$ is described by an equilibrium density $\pi := \rho_\infty^N(x) : N \to \mathbb{R}$. Then the Fokker-Planck equation is given by

$$\partial_t \rho = \text{div}_N \cdot \left( \pi \nabla_N \left( \frac{\rho}{\pi} \right) \right),$$

where $\nabla_N := \sum_{i=1}^d \tau_i^N \cdot \tau_i^N$ is surface gradient, $\nabla_i^N = \tau_i^N \cdot \nabla$ is the tangential derivative in the direction of $\tau_i^N$ and $\text{div}_N$ is the surface divergence defined as $\text{div}_N \xi = \sum_{i=1}^d \tau_i^N \cdot \nabla \tau_i^N \xi$.

4.1. **Construction of Voronoi tessellation and the upwind scheme on manifold $N$.** In this section, we construct an upwind scheme based on Voronoi tessellation for manifold $N$, which automatically gives a positive-preserving upwind scheme for the Fokker-Planck equation (4.1).

Suppose $(N, d_N)$ is a $d$ dimensional smooth closed submanifold of $\mathbb{R}^\ell$ and $d_N$ is induced by the Euclidean metric in $\mathbb{R}^\ell$. Let $Q := \{y_i\}_{i=1}^n$ be a set of cloud points well-distributed on $N$. Define the Voronoi cell as

$$C_i := \{ y \in N : d_N(y, y_i) \leq d_N(y, y_j) \text{ for all } y_j \in Q \},$$
with the volume $|C_i| = \mathcal{H}^d(C_i)$. Then $\mathcal{N} = \bigcup_{i=1}^n C_i$ is a Voronoi tessellation of manifold $\mathcal{N}$. One can see each $C_i$ is star shaped. Denote the Voronoi face for cell $C_i$ as
\begin{equation}
\Gamma_{ij} := C_i \cap C_j, \text{ and its area as } |\Gamma_{ij}| = \mathcal{H}^{d-1}(\Gamma_{ij})
\end{equation}
for any $j = 1, \cdots, n$. If $\Gamma_{ij} = \emptyset$ or $i \neq j$ then we set $|\Gamma_{ij}| = 0$.

Let $\chi_{C_i}$ be the characteristic function such that $\chi_{C_i} = 1$ for $y \in C_i$ and 0 otherwise. Then
\begin{equation}
\rho_{\text{approx}}(y) := \sum_{i=1}^n \rho_i \chi_{C_i}(y)
\end{equation}
is the piecewise constant probability distribution on $\mathcal{N}$ provided $\sum_{i=1}^n \rho_i |C_i| = 1$ and $\rho_i \geq 0$. Let $\pi_i$ be the approximated equilibrium density at $y_i$ satisfying $\sum_{i=1}^n \pi_i |C_i| = 1$. If $\rho_{\text{approx}}(y) = \sum_i \rho_i \chi_{C_i}(y)$ is an approximation of density $\rho_N(y)$, then $\rho_i$ is an approximation of the density $\rho_N(y_i)$.

Define the associated adjacent grids as
\begin{equation}
\text{VF}(i) := \{ j; \Gamma_{ij} \neq \emptyset \}.
\end{equation}
Then using the finite volume method and the divergence theorem on manifold, we have
\begin{equation}
\frac{d}{dt} \rho_i |C_i| = \frac{d}{dt} \int_{C_i} \rho_{\text{approx}} \mathcal{H}^d(C_i) = \sum_{j \in \text{VF}(i)} \int_{\Gamma_{ij}} \pi_i \nabla \mathcal{H}^{d-1}(\Gamma_{ij}) \cdot \mathbf{n},
\end{equation}
where $\mathbf{n}$ is the unit outward normal vector field on $\partial C_i$. Based on this, we introduce the following upwind scheme. For $i = 1, \cdots, n$,
\begin{equation}
\frac{d}{dt} \rho_i |C_i| = \frac{1}{2} \sum_{j \in \text{VF}(i)} \frac{\pi_i |\Gamma_{ij}|}{|y_i - y_j|} \left( \frac{\rho_j}{\pi_j} - \frac{\rho_i}{\pi_i} \right).
\end{equation}

We now interpret the upwind scheme as the forward equation for a Markov process with transition probability $P_{ji}$ (from $j$ to $i$) and jump rate $\lambda_j$
\begin{equation}
\frac{d}{dt} \rho_i |C_i| = \sum_{j \in \text{VF}(i)} \lambda_j P_{ji} \rho_j |C_j| - \lambda_i \rho_i |C_i|, \quad i = 1, 2, \cdots, n;
\end{equation}
where
\begin{equation}
\lambda_i := \frac{1}{|C_i| \pi_i} \sum_{j \in \text{VF}(i)} \frac{\pi_i |\Gamma_{ij}|}{|y_i - y_j|}, \quad i = 1, 2, \cdots, n;
\end{equation}
\begin{equation}
P_{ji} := \frac{1}{\lambda_j} \frac{\pi_i + \pi_j}{2 |C_j| \pi_j} |\Gamma_{ij}|, \quad j \in \text{VF}(i); \quad P_{ji} = 0, \quad j \notin \text{VF}(i).
\end{equation}
Assume $\pi_i > 0$ for all $i$, then we have $\lambda_i > 0$ for all $i$. One can see it satisfies $\sum_i P_{ji} = 1$ and the detailed balance property
\begin{equation}
P_{ji} \lambda_j \pi_j |C_j| = P_{ij} \lambda_i \pi_i |C_i|.
\end{equation}
We refer to [5] for the ergodicity of this Markov process.

In practice, instead of the $|C_i|, \Gamma_{ij}$ in (4.7), one shall use the approximated coefficients $\tilde{C}_i$ and $\tilde{\Gamma}_{ij}$ because we do not know the exact metric information of the manifold based only on point clouds. We omit the algorithm of finding the approximated $\tilde{C}_i$ and $\tilde{\Gamma}_{ij}$ and refer to [5, 4].
4.2. Unconditional stable explicit time stepping and exponential convergence. Now we propose an unconditionally stable explicit time discretization for the upwind scheme (4.6), which enjoys several good properties as the scheme (3.6), such as maximal principle, mass conservation law and exponential convergence.

Let \( \rho_i^k \) be the discrete density at discrete time \( k\Delta t \). To achieve both stability and efficiency, we introduce the following unconditional stable explicit time discretization for (4.7)

\[
\rho_i^{k+1} = \rho_i^k - \lambda_i \Delta t \rho_j^k + \Delta t \sum_{j \in V_F(i)} \lambda_i P_{ij} \rho_j^k, \quad i = 1, 2, \cdots, n
\]

which is

\[
\rho_i^{k+1} = \rho_i^k + \frac{\lambda_i \Delta t}{1 + \lambda_i \Delta t} \left( \sum_{j \in V_F(i)} P_{ij} \rho_j^k - \rho_i^k \right).
\]

For \( u_i^{k+1} := \frac{\rho_i^{k+1}}{\pi_i} \), the matrix formulation of (4.11) is

\[
u^{k+1} = (I + \Delta t \hat{B}) u^k,
\]

where

\[
\hat{B} := \{ \hat{b}_{ij} \} = \begin{cases} 
\frac{-\lambda_i}{1 + \lambda_i \Delta t}, & j = i; \\
\frac{\lambda_i P_{ij}}{1 + \lambda_i \Delta t}, & j \neq i
\end{cases} \quad \text{with} \quad \sum_j \hat{b}_{ij} = 0.
\]

We give the following proposition for several properties of scheme (4.10). The proof of this proposition is similar to Proposition 1 so we omit it.

**Proposition 3.** Let \( \Delta t \) be the time step and consider the explicit scheme (4.10).

Assume the initial data satisfies

\[
\sum_i (1 + \lambda_i \Delta t) \rho_i^0 |C_i| = \sum_i (1 + \lambda_i \Delta t) \pi_i |C_i|.
\]

Then we have

(i) the conversational law for \( g_i^{k+1} := (1 + \Delta t \lambda_i) \rho_i^{k+1} |C_i| \), i.e.

\[
\sum_i (1 + \lambda_i \Delta t) \rho_i^{k+1} |C_i| = \sum_i (1 + \lambda_i \Delta t) \rho_i^k |C_i|;
\]

(ii) the unconditional maximal principle for \( \frac{\rho_i^{k+1}}{\pi_i} \)

\[
\max_j \frac{\rho_j^{k+1}}{\pi_j} \leq \max_j \frac{\rho_j^k}{\pi_j};
\]

(iii) the \( \ell^\infty \) contraction

\[
\max_i \left| \frac{\rho_i^{k+1}}{\pi_i} - 1 \right| \leq \max_i \left| \frac{\rho_i^k}{\pi_i} - 1 \right|;
\]

(iv) the exponential convergence

\[
\left\| \frac{\rho_i^k}{\pi_i} - 1 \right\|_{\ell^\infty} \leq c |\mu_2|^k, \quad |\mu_2| < 1,
\]

where \( \mu_2 \) is the second eigenvalue of \( I + \Delta t \hat{B} \) (in terms of magnitude), i.e. \( \mu_2 = 1 - \text{gap}_B \Delta t \) and \( \text{gap}_B \) is the spectral gap of \( \hat{B} \).
4.3. Thresholding for shape dynamics. Assume the initial density is adjusted based on the mass conservation law (4.14). We now give the shape dynamics by combining the Fokker-Planck equation on manifold with the thresholding scheme. Assume initial data $\rho^0_i \in \{\rho^0_s, \rho^0_b\}$, which takes alternatively the value $\rho^0_s, \rho^0_b$. Assume the equilibrium is $\pi_i \in \{\pi_s, \pi_b\}$ which takes alternatively the value $\pi_s, \pi_b$.

Similar to Section 3.1, we choose the threshold constant $\xi_k$ at each step to conserve (4.14) as follows.

**Step 1.** Given $\rho^k_i \in \{\pi_s, \pi_b\}$, compute the explicit scheme (4.10) to update $\tilde{\rho}^{k+1}_i$ for any $i = 1, 2, \cdots, n$.

**Step 2.** Choose threshold constant $\xi^{k+1}$ and define

$$
\rho^{k+1}_i := \pi_s \chi\{i; \rho^{k+1}_i \leq \xi^{k+1}\} + \pi_b \chi\{i; \rho^{k+1}_i > \xi^{k+1}\}, \quad i = 1, 2, \cdots, n
$$

such that $\rho^{k+1}$ satisfies (4.14). Here $\xi^{k+1}$ can be found using bisection such that

$$
f(\xi^{k+1}) := \sum_i (1 + \lambda_i \Delta t) \rho^{k+1}_i |C_i| - \sum_i (1 + \lambda_i \Delta t) \pi_i |C_i| = 0.
$$

5. Computations. In this section, three numerical examples are carried out to examine the capability and efficiency of the equilibrium-driven deformation algorithm (EDDA), which are the RGB colored facial aging transformation, the pneumonia of COVID-19 invading and fading away on CT scan images and the continental evolution process.

5.1. Example I: RGB colored facial aging transformation. In this example, we have two images with the same size in the RGB color model showing a lady’s face at two different ages, and employ the model to simulate the transformation from one image (initial) to another image (equilibrium), which will illustrate the facial aging process with time. The strategy is to define each image as three matrices, each matrix containing values of a color mode (R or G or B). Then the transformation between the two images is computed by applying the inbetweening auto-animation three times based on the Fokker-Planck dynamics (2.1).

The two images are extracted from [7] and are both 355 pixels in width and 575 pixels in height, which means a total of 204125 pixel points in each image. The initial image data is first adjusted to meet the mass conservation law (3.11). Time step $\Delta t$ is set to 0.01 and the total number of iterations is set to 10000 thus the final iteration time $T = 100$. The horizontal resolution $\Delta x$ and $\Delta y$ are both $10^{-4}$. We use the unconditional stable explicit time stepping scheme (3.6) and the no-flux boundary condition (3.5) to the Fokker-Planck equation (2.1) in domain $\Omega$.

The relative root mean square errors (3.16) for the three color-modes are illustrated in Fig. 1 separately in semiology plot. Except for the different descend rates for the three colors, all simulated errors have the exponential convergence rates, which is consistent with the analysis in Proposition 1.

In order to see the transformation process between the two images, the images after iteration step 40, 100, 200, 400, 1000, 2000, 4000 and 10000 are shown and compared with the initial and the equilibrium images in Fig. 2. The transformation process between two images are fast in the beginning (e.g. before step 200) and relatively slow after then. The transformation process in Fig. 2 clearly reveals the potential changes in different parts of the lady’s face and hair with time. After 10000 steps of iterations, the updated image is nearly the same with the equilibrium except for the hair color.
5.2. Example II: COVID-19 pneumonia invading and fading away on CT scan images. In this section we will focus on an example based on the COVID-19 pneumonia invading and fading away process in a patient’s lung reflected on CT scan images and try to show the possible COVID-19 pneumonia growth dynamics with time before and after the treatment. In order to fulfill the task, two parts of simulations are presented. In the first part, two CT scan images taken on a patient’s lungs at the beginning (January 23th) and severe state (February 2nd) of the disease [6] are selected to be the initial and equilibrium state, respectively; see Fig. 3 (left). In the second part, two scan images at the severe state (February 2nd) and after a few-days’ treatment (February 15th) are selected to be the initial and equilibrium state, respectively; see Fig. 3 (right). Each CT scan image can be represented by a gray scale image matrix thus the same method in Example I can
Figure 3. Chest CT images of the critically severe COVID-19 patient [6]. The left column of figures illustrates the evading of pneumonia from January 23th to February 2nd and the right column illustrates the fading away of pneumonia from February 2nd to February 15th after treatment. Red circles indicate the significant COVID-19 pneumonia invading areas and blue circles indicate the significant pneumonia fading away areas.

be applied. The CT scan images are all cropped to 461 pixels in width and 370 pixels in height, which means a total of 170570 pixel points in each image. The time step $\Delta t$ is 0.01 and the total number of iterations is 6000 thus the final iteration time $T = 60$. The resolutions are $\Delta x = \Delta y = 10^{-4}$.

After 6000 iterations, the relative root mean square errors from two parts of simulations both decrease with an exponential rate, as is shown in Fig. 4. Moreover, the image evolution after step 20, 50, 100, 200, 500, 1000, 5000, 10000 (see Fig. 5) clearly demonstrate the pneumonia invading process into the patient’s lungs caused by COVID-19 in a few days (upper group of figures in Fig. 5) and the pneumonia fading away from the lungs after a stem cell treatment [6] is applied to the patient (lower group of figures in Fig. 5), indicating a potential success of this treatment [6]. We further compare the simulated pneumonia invading and fading process with the real CT scan images taken on January 30th and February 9th, respectively (see Fig. 3). The red circles in Fig. 3 clearly show the four significant COVID-19 pneumonia invading areas in the patient’s lungs, where the ground-glass opacity appears, showing damages to the lungs in the middle of the pneumonia invading process. As a comparison, the simulation results after Step 50 also show four areas where biggest density changes happen (red circles in Fig. 5) and they
Figure 4. The semilog plot of temporal evolution of the relative root mean square errors for COVID-19 pneumonia invading and fading away on CT scan images with the parameters $\Delta t = 0.01, T = 100$ and $\Delta x = \Delta y = 10^{-4}$. (up) The error evolution for the pneumonia invading process simulation. (down) The error evolution for the pneumonia fading away process simulation.

are consistent with the real invading areas. Similar comparisons can be carried out in the pneumonia fading process, in which the significant areas are indicated with blue circles in Fig. 3 and Fig. 5. The simulated results after Step 200 in (lower) Fig. 5 also show the two significant pneumonia fading areas in the patient’s lungs after treatment. The satisfactory agreements in this example indicate promising applications in this field.

5.3. Example III: Continental evolution process with thresholding for shape dynamics. In this section, we try to reveal the evolution process of continents in the world from Pangaea supercontinent (250 million years ago) to the current globe. In order to clearly distinguish the shape dynamics evolving the continents and the oceans, the thresholding scheme (4.19) and the Fokker-Planck dynamics (4.10) are combined to generate the inbetweening motions with the above sharp interfaces. The numerical experiment is carried out as follows.

**Step (i).** A group of points is selected on a unit sphere to be the dataset points. With the Centroidal Voronoi Tessellation (CVT) method on the unit sphere [8, 2], the Voronoi cells on the unit sphere are generated and the locations of dataset points are adjusted accordingly to ensure the uniformity of these cells. Thus, the distributions of continents and oceans derived from Pangaea period and current globe’s topography are described by two values (i.e. $\pi_s$ and $\pi_b$) on the Voronoi cells and are set to be the initial and equilibrium states, respectively. The Voronoi cell area $C_i, i = 1, \cdots, n$, with the total number of dataset points $n$, is computed and the Voronoi face $\Gamma_{ij}$ is determined by the geodesic length of the neighboring arc between cell $i$ and $j$.

**Step (ii).** Update the density at each point using the explicit scheme (4.10) for the linear Fokker-Planck equation.
The simulated COVID-19 pneumonia invading and fading away process on CT scan images with the parameters $\Delta t = 0.01$, $T = 100$ and $\Delta x = \Delta y = 10^{-4}$. Results after the step 20, 50, 100, 200, 500, 1000, 2000, 5000 are illustrated and compared to the initial and equilibrium scan images. The white part inside the lungs shown on images indicates the evidence of pneumonia. (up) The pneumonia invades into the patient’s lungs caused by COVID-19. (down) The pneumonia fades away from the lungs after a stem cell treatment is applied to the patient. Red circles indicate the significant COVID-19 pneumonia invading areas and blue circles indicate the significant pneumonia fading away areas.

**Step (iii).** After several linear iteration steps, the threshold is selected following the steps in Section 4.3 and the thresholding scheme is applied to update the data. The computations for Step (ii) and Step (iii) will be looped until reaches the total iteration steps. Besides, a simulation case which only evolves the linear Fokker-Planck equation is carried out as a comparison.

For example III, we select a total of 3000 dataset points and generate the Voronoi cells on the unit sphere via the CVT approximation algorithm; see Fig. 6. The standard deviation for all the cell areas is $3.2 \times 10^{-4}$, which means the nearly uniform distributions of data points on the sphere. The values at continental cells and the ocean cells are set to 0.9 and 0.1, respectively. The time step $\Delta t$ is set to 0.05 and
Figure 6. The unit sphere and the Voronoi cells on it. There are totally 3000 cells on the sphere. The black dots indicate locations of the point clouds on the sphere. The polygons with black edges are the Voronoi polygons generated with CVT algorithm.

Figure 7. The semilog plot of temporal evolution of relative root mean square errors for the continental evolution process with thresholding for shape dynamics with $\Delta t = 0.05$ and the total number of the thresholding adjustments is $N_t = 50$. The linear iterations before the $(k+1)$th thresholding adjustment is $2^k$. The red line indicates the error of simulations with only the linear Fokker-Planck algorithm while the blue line is the error of the linear algorithm combined with the thresholding scheme. The black dotted lines indicate the time steps when the thresholding adjustments are applied.

<table>
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Table 1. Comparison of the time steps needed to reach the tolerance between linear method and threshold method.

The total number of linear iterations before the $(k+1)$-th thresholding adjustment is set to $2^k$, $k = 1, \ldots, N_t$, where $N_t = 50$ is the times of the thresholding adjustments. The threshold $\xi_k$ is determined by a bisection method such that (4.20) is satisfied. Here, the bisection domain limitation criterion is set to $10^{-6}$. The total number of iterations for the comparison simulation is set to be 10000. We remark thresholding
The evolutions of continental movements on the unit sphere with the parameter $\Delta t = 0.05$ and the total number of the thresholding adjustments is $N_t = 50$. The continental evolution on the sphere after the 2th, 5th, 10th, 30th thresholding adjustment are illustrated and compared with the initial and equilibrium states. The black dots and polygons in each subplot illustrate the point clouds and the Voronoi cells, respectively. The orange and blue patch indicate the land and ocean, respectively. ‘TH’ is short for ‘thresholding step’. The formation of the Antarctic is revealed at the bottom (southern part) of the globe (black arrow in TH 5). Note that the globes are shown in the same view angle so the Antarctic continental is out of view in the last two subplots.

adjustment after each $2k$, $k = 1, 2, \cdots, N_t$ Fokker-Planck iterations is different from the one with fixed-time step thresholding adjustment whose continuum limit may lead to a motion by mean curvature correction. In our scheme, as $k$ increases, the leading contribution still comes from the linear Fokker-Planck dynamics. So analogue to reinitialization in the level-set method, our thresholding adjustment can be regarded as a reinitialization at each $(k+1)$-th adjustment step, which is an efficient numerical method for the shape dynamics simulation and has finite time convergence property to its equilibrium; see Proposition 2.

Fig. 7 shows the temporal variations of the relative root mean square errors for the numerical example in the first 1200 time steps. The error from the thresholding method generally have a descend trend although with some abrupt increase due to the thresholding adjustments. The error decreases to nearly zero (less than the machine accuracy) after the 30th thresholding adjustment (a total of 960 time steps), which indicates the data is updated to the equilibrium. As a comparison,
the error of the simulation via the linear method (red line in Fig. 7), which leads an exponential convergence rate, is smaller than that from thresholding method before the 960th time step (black circle in Fig. 7) and is larger after then. In order to further compare the efficiency of the two methods, we calculated the total time steps needed for the error to reach difference tolerances and listed them in Table 1. Based on the finite time convergence property in Proposition 2, as long as the tolerance between the Fokker-Planck solution and the equilibrium is smaller than a specific quantity, then after the thresholding adjustment, the resulting binary density will stay at equilibrium. Thus the fixed number of time steps in Table 1 independent of the tolerance of error is consistent with Proposition 2. The comparisons clearly reveal the efficiency of the thresholding adjustment in the application of the shape dynamics, especially when the tolerance is small.

The continental evolution on the sphere after the 2th, 5th, 10th, 30th thresholding adjustment is illustrated and compared with the initial and equilibrium states in Fig. 8. After several steps of thresholding adjustment, the sharp shapes of continents quickly move from the initial Pangaea supercontinent towards the equilibrium state of current continents. The distributions of continents and oceans reaches the equilibrium state after the 28th thresholding adjustment, exactly the same as the current distributions. Although the evolution of the continental movements is simulated with the data-driven model, some potential dynamics of continental drifting such as the Antarctic formation can be noticed in the evolutions in Fig. 8, which may contribute to the detailed explanation of the continental drifting theory.

6. Discussion. We propose an efficient and universal equilibrium-driven deformation algorithm (EDDA) to simulate the inbetweening transformations given an initial and equilibrium. The algorithm automatically cooperates positivity, unconditional stability, mass conservation law, exponentially convergence and also the manifold structure suggested by dataset. Using EDDA, three challenging examples, (I) facial aging process, (II) COVID-19 pneumonia invading/fading process, and (III) continental evolution process are conducted efficiently. EDDA is shown to be a very efficient and universal method with enormous potential applications in other fields of science and industry.

Acknowledgments. The authors would like to thank Prof. Haiyan Gao for helpful suggestions.

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Received May 2020; revised October 2020.

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