Perron-Frobenius operator filter for stochastic dynamical systems

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Abstract

The filtering problems are derived from a sequential minimization of a quadratic function representing a compromise between model and data. In this paper, we use the Perron-Frobenius operator in stochastic process to develop a Perron-Frobenius operator filter. The proposed method belongs to Bayesian filtering and works for non-Gaussian distributions for nonlinear stochastic dynamical systems. The recursion of the filtering can be characterized by the composition of Perron-Frobenius operator and likelihood operator. This gives a significant connection between the Perron-Frobenius operator and Bayesian filtering. We numerically fulfil the recursion through approximating the Perron-Frobenius operator by Ulam's method. In this way, the posterior measure is represented by a convex combination of the indicator functions in Ulam's method. To get a low rank approximation for the Perron-Frobenius operator filter, we take a spectral decomposition for the posterior measure by using the eigenfunctions of the discretized Perron-Frobenius operator. A convergence analysis is carried out and shows that the Perron-Frobenius operator filter achieves a higher convergence rate than the particle filter, which uses Dirac measures for the posterior. The proposed method is explored for the data assimilation of the stochastic dynamical systems. A few numerical examples are presented to illustrate the advantage of the Perron-Frobenius operator filter over particle filter and extend Kalman filter.

keywords: Perron-Frobenius operator, Bayesian filtering, stochastic dynamical systems, particle filter

1 Introduction

In recent years, the operator-based approach has been extensively exploited to analyze dynamical systems. The two primary candidates of the approach are Perron-Frobenius operator and its dual operator, Koopman operator. Many data-driven methods have been developed for numerical approximation of these operators. The two operators are motivated to approximate the dynamical system's behavior from different perspectives. The Koopman operator is used to study the evolution of observations, while Perron-Frobenius operator (PFO) characterizes the transition of densities. Therefore, the PFO deals with the system's uncertainties

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in the form of probability density functions of the state. In practice, it determines an absolutely continuous probability measure preserved by a given measurable transformation on a measure space.

The Perron-Frobenius operator has been widely used to characterize the global asymptotic behavior of dynamical systems derived from many different domains such as fluid dynamics [1], molecular dynamics [2], meteorology and atmospheric sciences [3], and to estimate invariant sets or metastable sets with a toolbox like in [4]. It is of great interest to study the invariant density of Perron-Frobenius operator [5] and design efficient numerical approaches. Then one can apply ergodic theorems to the statistical properties of deterministic dynamical systems.

Since PFO is able to transport density of a Markov process, its approximation is necessary for numerical model transition probability of the Markov process. Many different numerical methods [6], such as Ulam's method and Petrov-Galerkin method, are proposed for approximation of the Perron-Frobenius operator. As the PFO operates on infinite-dimensional spaces, it is natural to project it onto a finite-dimensional subspace spanned by suitable basis functions. The projection is usually accomplished by Galerkin methods with weak approximation. It was originally proposed by Ulam [7], who suggested that one can study the discrete Perron-Frobenius operator on the finite-dimensional subspace L^1 of indicator functions according to a finite partition of the region. Convergence analysis of Ulam's method is discussed in many literatures [8, 9].

The classical filtering problems in stochastic processes are investigated in [10, 11]. In the paper, the models of filtering problems are considered with discrete-time and continuoustime stochastic processes defined by the solutions of SDEs, which can model a majority of stochastic dynamical systems in the real world. The filtering methods have been widely used for geophysical applications, such as oceanography [12], oil recovery [13], atmospheric science, and weather forecast [14]. Remarkably, as one of the filtering methods, Kalman filter [15] has been well-known for low-dimensional engineering applications in linear Gaussian models, and it has been also developed and utilized in many other fields [16–18]. For nonlinear problems, the classical filters, such as 3DVAR [19], Extended Kalman filter [10] and Ensemble Kalman filter [20], usually invoke a Gaussian ansatz. They are often used in the scenarios with small noisy observation and high dimensional spaces. However, these extensions rely on the invocations of Gaussian assumption. As a sequential Monte Carlo method, particle filter [21] is able to work well for the nonlinear and non-Gaussian filtering problems. It can be proved to estimate true posterior filtering problems in the limit of large number of particles.

Although the particle filter (PF) can treat the nonlinear and non-Gaussian filtering problems, it has some limitations in practice. First of all, particle filter handles well in lowdimensional systems, but it may occur degeneracy [22] when the systems have very large scale. It means that the maximum of the weights associated with the sample ensemble converges to one as the system dimension tends to infinity. To avoid degeneracy, it requires a great number of particles that scales exponentially with the system dimension. This is a manifestation of the curse of dimensionality. Resampling, adding jitter and localisation are introduced to circumvent this curse of dimensionality and get the accurate estimation of highdimensional probability density functions (PDFs). The particle filter also does not perform well in geophysical applications of data assimilation, because the data in these application have strongly constraints on particle location [23]. This impacts on the filtering performance. Besides, the prior knowledge of the transition probability density functions in particle filter is necessary to be known, and the efficient sampling methods such as acceptance-rejection method and Markov chain Monte Carlo, need to be used for complicated density functions. The sampling is particularly a challenge in high dimensional spaces. To overcome these difficulties, we propose a Perron-Frobenius operator filter (PFOF), which does not use particles and any sampling methods. The information of prior probability density is not required in the method, which needs data information instead. The method works well in nonlinear and non-Gaussian models.

In this paper, we propose PFOF to treat nonlinear stochastic filtering problems. The method transfer filtering distribution with the Perron-Frobenius operator. For filtering problems, the update of filtering distribution involves two steps: predication and analysis. In prediction, the density is transported with a transition density function given by a Markov chain of the underlying system. In analysis, the density is corrected by Bayes' rule under the likelihood function given by observations. Thus, the update of filtering distribution can be expressed as a composition of PFO and likelihood functions. In the simulation process, Ulam' method is used to discretize the PFO and project it onto a finite-dimensional space of indicator functions. Hence the filtering density is also projected onto the subspace and is represented by a linear convex combination of the indicator functions. The recursion of filtering distribution is then expressed by a linear map of weights vectors associated with the basis functions via the PFO and likelihood function. For the high dimensional problems, we propose a low-rank PFOF (lr-PFOF) using a spectral decomposition. To this end, we first use the eigenfunctions of the discretized PFO to represent the spectral decomposition of the density functions. Then we make a truncation of the decomposition and use the eigenfunctions corresponding to the first few dominant eigenvalues. This can improve the online assimilation efficiency. The idea of PFO is extended to the continuous-time filtering problems. In these problems, Zakai equation characterizes the transition of filtering density. We utilize the approximation of the Perron-Frobenius operator to compute the Zaikai equation and obtain the posterior density functions of the continuous-time filtering problems.

We compare the proposed method with the particle filter. For PFOF, we give an error estimate in the total-variance distance between the approximated posterior measure and the truth posterior. The estimate implies that PFOF achieves a convergence rate $O(\frac{1}{N})$, which is faster than particle filters with the same number N of basis functions. Our numerical results show that PFOF also renders better accuracy than that of extended Kalman filter.

The rest of the paper is organized as follows. In Section 2, we express the Bayesian filter in terms of the Perron-Frobenius operator. In Section 3, we present the recursion of the filtering empirical measure with an approximated Perron-Frobenius operator. Then we derive PFOF as well as lr-PFOF, and analyze an error estimate subsequently. PFOF is also extended to the continuous-time filtering problems. A comprehensive comparison with particle filter is give in Section 4. A few numerical results of stochastic filtering problems are given in Section 5. Section 6 concludes the paper in a summary.

2 Preliminaries

We give a background review on Perron-Frobenius operator [24] (PFO) and Bayesian filter in this section. The Perron-Frobenius operator transports the distributions over state space and describes the stochastic behavior of the dynamical systems. The framework of Bayesian filter is introduced and summarized as a recursive formula with PFO.

2.1 Perron-Frobenius operator

Let X be a metric space, \mathcal{B} the Borel- σ -algebra on X, and $\Phi : X \to X$ a nonsingular transformation. Let \mathcal{M} denote the space of all finite measures on (X, \mathcal{B}) and μ is a finite measure. The phase space is defined on a measure space (X, \mathcal{B}, μ) . The Perron-Frobenius operator $\mathcal{P} : \mathcal{M} \to \mathcal{M}$ is a linear and infinite-dimensional operator defined by

$$\mathcal{P}\mu(A) = \mu(\Phi^{-1}(A)), \quad \forall A \in \mathcal{B}.$$
 (2.1)

The PFO is a linear, positive and non-expansive operator, and hence a Markov operator. We can also track the action on distributions in the function space. In the paper, we denote $L^1(X) := L^1(X, \mathcal{B}, \mu)$. Let $f \in L^1(X)$ be the probability density function (PDF) of a X-valued random variable x. Since Φ is a nonsingular with respect to μ , there is a $g \in L^1(X)$ satisfying $\int_{\Phi^{-1}(A)} f d\mu = \int_A g d\mu$ for all $A \in \mathcal{B}$. Then g is the function characterizing the distribution of $\Phi(x)$. The mapping $f \mapsto g$, defined uniquely by a linear operator $\mathcal{P} : L^1(X) \to L^1(X)$:

$$\int_{A} \mathcal{P}f \, d\mu = \int_{\Phi^{-1}(A)} f \, d\mu, \quad \forall A \in \mathcal{B}.$$
(2.2)

The operator \mathcal{P} is called the Perron-Frobenius operator. With the definition (2.1) and (2.2), we make the connection between probability density function and the measure associated with the PFO. When f is a probability density function with respect to an absolutely continuous probability measure $\mu \in \mathcal{M}(X)$, g is another PDF with respect to the absolutely continuous probability measure $\mu \circ \Phi^{-1}$. In addition, the measure $\mu \in \mathcal{M}(X)$ is an invariant measure of \mathcal{P} when $\mathcal{P}\mu = \mu$ holds.

Let $\Psi : \mathbb{R}_+ \times X \to X$ be a nonsingular flow map for a deterministic continuous-time dynamical system. Then $\Psi_\tau : X \to X$ is nonsingular for each $\tau \in \mathbb{R}^+$. The transfer operator $\mathcal{P}_\tau : L^1(X) \to L^1(X)$ is time-dependent and has an analogous definition to (2.2), such that

$$\int_{A} \mathcal{P}_{\tau} f \, d\mu = \int_{\Psi_{\tau}^{-1}(A)} f \, d\mu$$

The $\{\mathcal{P}_{\tau}\}_{\tau\geq 0}$ forms a semigroup of the Perron-Frobenius operators. We note that $\{\mathcal{P}_{\tau}\}_{\tau\geq 0}$ has an infinitesimal generator \mathscr{A}_{PF} by Hille-Yosida Theorem.

Let us consider the Perron-Frobenius operator in stochastic dynamic systems and the infinitesimal generator of PFO associated to the stochastic solution semiflow induced by a stochastic dynamical equation (SDE). Let $b : X \to X$ and $\sigma : X \to X$ be smooth time-invariant functions. Suppose that a stochastic process x_t is the solution to the time-homogeneous stochastic differential equation:

$$dx_t = b(x_t)dt + \sigma(x_t)dW_t, \quad x(t_0) \sim \rho_0, \tag{2.3}$$

where W_t is a standard Brownian motion. In this case, the distribution of the stochastic process x_t can be described by a semigroup of Perron-Frobenius operators $\{\mathcal{P}_{\tau}\}_{\tau>0}$ on $L^1(X)$. The generator of $\{\mathcal{P}_{\tau}\}_{\tau>0}$ is a second-order differential operator on X. The PDE defined by the generator describes the evolution of the probability density of x_t .

Suppose that Φ is the mapping of the stochastic dynamical system (2.3) and $\Phi(x)$ is an *X*-valued random variable over the probability space (X, \mathcal{B} , μ). Given a stochastic transition function $p_{\tau} : X \times \mathcal{B} \to [0, 1]$ induced by Φ , we consider probability measure μ translated with a linear operator defined in terms of the transition function $p_{\tau}(x, \cdot)$. The stochastic PFO [25] $\mathcal{P}_{\tau} : \mathcal{M} \to \mathcal{M}$ is defined by

$$\mathcal{P}_{\tau}\mu(A) = \int_{X} p_{\tau}(x, A) \, d\mu(x), \quad \forall A \in \mathcal{B}.$$
(2.4)

If $p_{\tau}(x, \cdot)$ is absolutely continuous to μ for all $x \in X$, there exists a nonnegative transition density function $q_{\tau}: X \times X \to \mathbb{R}$ with $q_{\tau}(x, \cdot) \in L^{1}(X)$ and

$$P(x_{t+\tau} \in A | x_t = x) = \int_A q_\tau(x, y) d\mu(y), \quad \forall A \in \mathcal{B}.$$

The transition density function is the infinite-dimensional counterpart of the transition matrix for a Markov chain. Now we define the stochastic PFO associated with transition density. If $f \in L^1(X)$ is a probability density function, the Perron-Frobenius semigroup of operators $\mathcal{P}_{\tau}: L^1(X) \to L^1(X), \tau > 0$, is defined by

$$\mathcal{P}_{\tau}f(y) = \int_X q_{\tau}(x,y)f(x) \, d\mu(x).$$

The PFO \mathcal{P}_{τ} defined here translates the probability density function of x_t with time. Let ρ be a probability density. The infinitesimal generator \mathscr{A}_{PF} of \mathcal{P}_{τ} is given by

$$\mathscr{A}_{PF}\rho = -\nabla \cdot (b\rho) + \frac{1}{2}\nabla \cdot \nabla \cdot (\sigma\sigma^{T}\rho).$$

We assume that $\tilde{\rho}: [0, \infty) \times X \to [0, \infty)$ is the probability density function of the solution x_t in (2.3) and ρ_0 is the density function of the initial condition x_0 . Then $\tilde{\rho}$ solves the Fokker-Planck equation,

$$\begin{cases} \frac{\partial \widetilde{\rho}}{\partial t} = -\nabla \cdot (b\widetilde{\rho}) + \frac{1}{2}\nabla \cdot \nabla \cdot (\sigma\sigma^T \widetilde{\rho}), \quad (t, x) \in (0, \infty) \times X, \\ \widetilde{\rho}(0, x) = \rho_0(x). \end{cases}$$

If the phase space X is compact and $b \in C^3(X, X)$, the equation has a unique solution, which is given by

$$\widetilde{\rho}(t,x) = \mathcal{P}_t \rho_0(x).$$

2.2 Bayesian filter

In this section, we present the framework of Bayesian filter in discrete time from the perspective of Bayes' rule. In filtering problems, a state model and an observation model are combined to estimate the posterior distribution, which is a conditional distribution of the state given by observation. Let us consider the dynamical model governed by the flow $\Psi \in C(\mathbb{R}^n, \mathbb{R}^n)$ with noisy observations $y = \{y_j\}_{j \in \mathbb{Z}^+}$ depending on the function $h(x) : \mathbb{R}^n \to \mathbb{R}^p$:

$$\begin{cases} x_{j+1} = \Psi(x_j) + \xi_j, \ j \in \mathbb{N}, \ x_0 \sim \rho_0, \\ y_{j+1} = h(x_{j+1}) + \eta_{j+1}, \ j \in \mathbb{N}, \end{cases}$$
(2.5)

where $\xi := \{\xi_j\}_{j \in \mathbb{N}}$ is an i.i.d. sequence with $\xi_j \sim N(0, \Sigma)$ and $\eta := \{\eta_j\}_{j \in \mathbb{Z}^+}$ is an i.i.d. sequence with $\eta_j \sim N(0, R)$. Let $Y_j = \{y_l\}_{l=1}^j$ denote the data up to time t_j . The filtering problem aims to determine the posterior PDF $p(x_j|Y_j)$ of the random variable $x_j|Y_j$ and the sequential updating of the PDF as the data increases. The Bayesian filtering involves two steps: prediction and analysis. It provides a derivation of $p(x_{j+1}|Y_{j+1})$ from $p(x_j|Y_j)$. The prediction is concerned with the map $p(x_j|Y_j) \mapsto p(x_{j+1}|Y_j)$ and the analysis derives the map $p(x_{j+1}|Y_j) \mapsto p(x_{j+1}|Y_{j+1})$ by Bayes's formula.

Prediction

$$p(x_{j+1}|Y_j) = \int_{\mathbb{R}^n} p(x_{j+1}|Y_j, x_j) p(x_j|Y_j) dx_j$$

=
$$\int_{\mathbb{R}^n} p(x_{j+1}|x_j) p(x_j|Y_j) dx_j.$$
 (2.6)

Note that $p(x_{j+1}|Y_j, x_j) = p(x_{j+1}|x_j)$, because Y_j provides indirect information about determining x_{j+1} . Since $p(x_{j+1}|x_j)$ is specified by the underlying model (2.5) and

$$p(x_{j+1}|x_j) \propto \exp(-\frac{1}{2}|\Sigma^{-\frac{1}{2}}(x_{j+1} - \Psi(x_j))|^2),$$
 (2.7)

the prediction provides the map from $p(x_j|Y_j)$ to $p(x_{j+1}|Y_j)$. Let $\hat{\mu}_j$ be the prior probability measure corresponding to the density $p(x_j|Y_{j-1})$ and μ_j be the posterior probability measure on corresponding to the density $p(x_j|Y_j)$. The stochastic process $\{x_j, j \in \mathbb{N}\}$ of (2.5) is a Markov chain with the transition kernel $p(\cdot, \cdot)$ determined by $p(x_j, x_{j+1}) = p(x_{j+1}|x_j)$. Then we can rewrite (2.6) as

$$\widehat{\mu}_{j+1}(\cdot) = (\mathcal{P}\mu_j)(\cdot) := \int_{\mathbb{R}^n} p(x_j, \cdot)\mu_j(dx_j) = \int_{\mathbb{R}^n} p(x_j, \cdot)d\mu(x_j)^1.$$
(2.8)

In particular, the operator \mathcal{P} coincides with the Perron-Frobenius operator defined in (2.4).

Analysis

$$p(x_{j+1}|Y_{j+1}) = p(x_{j+1}|Y_j, y_{j+1})$$

$$= \frac{p(y_{j+1}|x_{j+1}, Y_j)p(x_{j+1}|Y_j)}{p(y_{j+1}|Y_j)}$$

$$= \frac{p(y_{j+1}|x_{j+1})p(x_{j+1}|Y_j)}{p(y_{j+1}|Y_j)}.$$
(2.9)

Note that $p(y_{j+1}|x_{j+1}, Y_j) = p(y_{j+1}|x_{j+1})$ and Bayes's formula is used in the second equality. The likelihood function $p(y_{j+1}|x_{j+1})$ is determined by the observation model: $p(y_{j+1}|x_{j+1}) \propto$

¹Refer to [26], if the function $f \in L^1(X)$ on a measure space (X, \mathcal{B}, μ) is said to be μ integrable, we have $\int f d\mu = \int f(x) \mu(dx) = \int f(x) d\mu(x)$.

 $\exp(-\frac{1}{2}|R^{-\frac{1}{2}}(y_{j+1}-h(x_{j+1}))|^2)$. Let

$$g_j(x_{j+1}) := \exp(-\frac{1}{2}|R^{-\frac{1}{2}}(y_{j+1} - h(x_{j+1}))|^2).$$
(2.10)

The analysis provides a map from $p(x_{j+1}|Y_j)$ to $p(x_{j+1}|Y_{j+1})$, so we can represent the update of the measure $\mu_{j+1}(\cdot)$ by

$$\mu_{j+1}(\cdot) = (L_j \widehat{\mu}_{j+1})(\cdot) := \frac{g_j(x_{j+1})\widehat{\mu}_{j+1}(\cdot)}{\int_{\mathbb{R}^n} g_j(x_{j+1})\widehat{\mu}_{j+1}(\cdot)},$$
(2.11)

where the likelihood operator L_j is defined by

$$(L_{j}\mu)(dx) = \frac{g_{j}(x)\mu(dx)}{\int_{\mathbb{R}^{n}} g_{j}(x)\mu(dx)}.$$
(2.12)

In general, the prediction and analysis provide the mapping from μ_j to μ_{j+1} . The prediction maps μ_j to $\hat{\mu}_{j+1}$ through the Perron-Frobenius operator \mathcal{P} , while the analysis maps $\hat{\mu}_{j+1}$ to μ_{j+1} through the likelihood operator L_j . Then we represent the μ_{j+1} using formulas (2.8) and (2.11), and summarize Bayesian filtering as

$$\mu_{j+1} = L_j \mathcal{P} \mu_j. \tag{2.13}$$

The μ_0 is assumed to be a known initial probability measure. We note that \mathcal{P} does not depend on j, because the prediction step is governed by the same Markov process at each j. However, L_j depends on j because the different observations are used to compute the likelihood at each j. In this way, the evolution of μ_j processes through a linear operator \mathcal{P} and a nonlinear operator L_j . The approximation of μ_j can be achieved by the numerical iteration of (2.13).

3 Bayesian filter in terms of Perron-Frobenius operator

It is noted that the Perron-Frobenius operator translates a probability density function with time according to the flow of the dynamics. We extend the idea to filtering problems to represent the transition of the posterior probability density function, i.e., the filtering distribution. Therefore, we propose a filtering method: Perron-Frobenius operator filter (PFOF). In the proposed method, the density function is projected onto an approximation subspace spanned by indicator functions. The fluctuation of the density function, which is approximated by weights vector, is transferred by PFO and likelihood operator. Moreover, we present a low-rank Perron-Frobenius operator filter (lr-PFOF), which is a modified version of the PFOF.

3.1 Perron-Frobenius operator filter

The iteration (2.13) is helpful to design a filter method. According to definition (2.8), the operator \mathcal{P} in the iteration is Perron-Frobenius operator corresponding to the flow Ψ of the

model (2.5). Based on the idea, we propose a Perron-Frobenius operator filter, which utilizes Ulam's method [7] to approximate operator \mathcal{P} in the iteration. In PFOF, we simply use \mathcal{P} for \mathcal{P}_{τ} because the discrete time steps of the state model keep the same. In this manner, the iteration of filtering distribution of PFOF becomes

$$\mu_{j+1}^N = L_j \mathcal{P}^N \mu_j^N, \quad \mu_0^N = \mu_0, \tag{3.14}$$

where \mathcal{P}^N calculated by the Ulam's method is an approximation of \mathcal{P} . Ulam's method is a Galerkin projection method to discretize the Perron-Frobenius operator. We first give a discretisation of the phase space. Let $B = \{\mathbb{B}_1, \dots, \mathbb{B}_N\} \subset \mathcal{B}$ be a finite number of measure boxes and a disjoint partition of phase space X. The indicator function is a piecewise constant function and is defined by

$$\mathbb{1}_{\mathbb{B}_i}(x) = \begin{cases}
1, & \text{if } x \in \mathbb{B}_i, \\
0, & \text{otherwise.}
\end{cases}$$
(3.15)

Ulam proposed to use the space of a family of indicator functions $\{\mathbb{1}_{\mathbb{B}_1}, \cdots, \mathbb{1}_{\mathbb{B}_N}\}$ as the approximation space for the PFO. We define the projection space $V_N := \text{span}\{\mathbb{1}_{\mathbb{B}_1}, \cdots, \mathbb{1}_{\mathbb{B}_N}\}$. The $V_N \in L^1(X)$ is regarded as an approximation subspace of $L^1(X)$. For each $\rho \ge 0$ in $L^1(X)$, we define the operator $\pi_N : L^1(X) \to V_N$ by

$$\pi_N \rho = \sum_{i=1}^N \omega^{(i)} \mathbb{1}_{\mathbb{B}_i}, \quad \text{where} \quad \omega^{(i)} := \frac{\int_{\mathbb{B}_i} \rho \, d\mu}{\mu(\mathbb{B}_i)}.$$
(3.16)

Then π_N is the projection onto V_N . Due to the projection, we define the discretized PFO \mathcal{P}^N as

$$\mathcal{P}^N = \pi_N \circ \mathcal{P}.$$

We can represent the linear map $\mathcal{P}^N|_{V_N^1}: V_N^1 \to V_N^1$, where $V_N^1:=\{f \in V_N: \int |f|d\mu=1\}$ by a matrix $P^N = (P_{ij}^N) \in \mathbb{R}^{N \times N}$ whose entries $P_{ij}^N = \frac{1}{\mu(\mathbb{B}_i)} \int_{\mathbb{B}_i} \mathcal{P}\mathbb{1}_{\mathbb{B}_j} d\mu$. The entries characterizes the transition probability from the box \mathbb{B}_i to box \mathbb{B}_j under the flow map Ψ . We can show

$$P_{ij}^N = \frac{\mu(\mathbb{B}_i \cap \Psi^{-1}(\mathbb{B}_j))}{\mu(\mathbb{B}_i)}.$$
(3.17)

A Markov chain for Ψ arises as the discretization \mathcal{P}^N of the PFO, and the Markov chain has transition matrix P^N . So the Ulam's method can be described either in terms of the operator \mathcal{P}^N or the matrix P^N . By the projection π_N , the density ρ can be expressed as a vector $\mathbf{W} = [\omega^{(1)}, \dots, \omega^{(N)}]$, where $\omega^{(i)}$ is the weight of the basis function $\mathbb{1}_{\mathbb{B}_i}$. Since the entries P_{ij}^N represent the transition probability from \mathbb{B}_i to \mathbb{B}_j , they can be estimated by a Monte-Carlo method, which gives a numerical realization of Ulam's method [6]. We randomly choose a large number of points $\{x_i^l\}_{l=1}^n$ in each \mathbb{B}_i and count the number of times $\Psi(x_i^l)$ contained in box \mathbb{B}_j . Then P_{ij}^N is calculated by

$$P_{ij}^N \approx P_{n,ij}^N = \frac{1}{n} \sum_{l=1}^n \mathbb{1}_{\mathbb{B}_j}(\Psi(x_i^l)).$$
 (3.18)

The Monte-Carlo method is used as an approximation to the integrals (3.17). The convergence of the Ulam's method depends on the choice of the partition of the region and the number of points. Based on indicator basis functions, we denote the the empirical density in the PFOF with respect to the measure μ_i^N as

$$\rho_j^N(x) = \sum_{i=1}^N \omega_j^{(i)} \mathbb{1}_{\mathbb{B}_i}(x), \qquad (3.19)$$

where $\mathbb{1}_{\mathbb{B}_i}(\cdot)$ is the indicator function defined by (3.15) and j represents the index of time sequence. In this way, the density ρ_j^N can be represented by the vector of the weights $\mathbf{W}_j = [\omega_j^{(1)}, \cdots, \omega_j^{(N)}]$. Suppose that $\mathbf{W}_j = [\omega_j^{(1)}, \cdots, \omega_j^{(N)}]$ and $\mathbf{W}_{j+1} = [\omega_{j+1}^{(1)}, \cdots, \omega_{j+1}^{(N)}]$ are separately the weights of $\pi_N \rho_j$ and $\pi_N \rho_{j+1} := \pi_N \mathcal{P} \rho_j$. When the region is evenly divided, the evolution of density functions becomes a Markov transition equation of the weights:

$$\mathbf{W}_{i+1} = \mathbf{W}_i P^N, \tag{3.20}$$

where P^N is the matrix form of discretized PFO. We consider the projection of the $\mathcal{P}\rho_j$, i.e.,

$$\pi_N \mathcal{P} \rho_j = \sum_{i=1}^N \omega_{j+1}^{(i)} \mathbb{1}_{\mathbb{B}_i}.$$
(3.21)

In addition, note that

$$\pi_N \mathcal{P} \rho_j = \pi_N \mathcal{P} \sum_{i=1}^N \omega_j^{(i)} \mathbb{1}_{\mathbb{B}_i} = \sum_{i=1}^N \omega_j^{(i)} \pi_N (\mathcal{P} \mathbb{1}_{\mathbb{B}_i}).$$

We denote $\pi_N(\mathcal{P}\mathbb{1}_{\mathbb{B}_i}) = \sum_{k=1}^N c_{ik}\mathbb{1}_{\mathbb{B}_k}$, where

$$c_{ik} = \frac{\int_X \mathcal{P}(\mathbb{1}_{\mathbb{B}_i})\mathbb{1}_{\mathbb{B}_k} dx}{\mu(\mathbb{B}_k)} = \frac{\int_{\mathbb{B}_k} \mathcal{P}(\mathbb{1}_{\mathbb{B}_i}) dx}{\mu(\mathbb{B}_k)}$$
$$= \frac{\int_{\Psi^{-1}(\mathbb{B}_k)} \mathbb{1}_{\mathbb{B}_i} dx}{\mu(\mathbb{B}_k)} = \frac{\mu(\mathbb{B}_i \cap \Psi^{-1}(\mathbb{B}_k))}{\mu(\mathbb{B}_k)}.$$

Then we have

$$\pi_N \mathcal{P} \rho_j = \sum_{i=1}^N \omega_j^{(i)} \sum_{k=1}^N c_{ik} \mathbb{1}_{\mathbb{B}_k} = \sum_{k=1}^N \sum_{i=1}^N \omega_j^{(i)} c_{ik} \mathbb{1}_{\mathbb{B}_k}$$
$$= \sum_{k=1}^N \sum_{i=1}^N \frac{\mu(\mathbb{B}_i)}{\mu(\mathbb{B}_k)} \omega_j^{(i)} P_{ik}^N \mathbb{1}_{\mathbb{B}_k}.$$

Comparing to (3.21), we get

$$\omega_{j+1}^{(k)} = \sum_{i=1}^{N} \frac{\mu(\mathbb{B}_i)}{\mu(\mathbb{B}_k)} \omega_j^{(i)} P_{ik}^N.$$
(3.22)

Thus, if we give a uniform partition of the X, i.e., $\mu(\mathbb{B}_i) = \mu(\mathbb{B}_j)$, $\forall i, j \in N$, then we get the result (3.20). With the expression, we design the following prediction step and analysis step to approximate the posterior distribution $p(x_{j+1}|Y_{j+1})$.

Prediction In this step, we give a set of boxes $\{\mathbb{B}_1, \dots, \mathbb{B}_N\} \subset \mathcal{B}$, which is a uniform partition of X, and denote the mass point of each box as $x^{(i)}, i = 1, \dots, N$. Define $\widehat{\mathbf{W}}_j = [\widehat{\omega}_j^{(1)}, \dots, \widehat{\omega}_j^{(N)}]$ as the prior weight vector and $\mathbf{W}_j = [\omega_j^{(1)}, \dots, \omega_j^{(N)}]$ as the posterior weight vector. In equation (2.8), we note that the prior density $p(x_{j+1}|Y_j)$ is computed under the linear operator \mathcal{P} . To discretize the formula $\widehat{\mu}_{j+1} = \mathcal{P}\mu_j$, we build a map between the weights of the density function,

$$\widehat{\mathbf{W}}_{j+1} = \mathbf{W}_j P^N.$$

The formula contains the prior information of the underlying system (2.5) because the PFO in the formula is defined by the transition kernel p of the system. With the basis functions $\mathbb{1}_{\mathbb{B}_i}(\cdot)$, the empirical prior measure is given by

$$\widehat{\mu}_{j+1}^N = \sum_{i=1}^N \widehat{\omega}_{j+1}^{(i)} \mathbb{1}_{\mathbb{B}_i}(dx).$$

Analysis In this step, we derive the posterior measure μ_{j+1}^N . To achieve this, we apply Bayes's formula (2.9) on weights and update the weights by

$$\omega_{j+1}^{(i)} = \widetilde{\omega}_{j+1}^{(i)} / (\sum_{n=1}^{N} \widetilde{\omega}_{j+1}^{(n)}), \quad \widetilde{\omega}_{j+1}^{(i)} = g_j(x^{(i)}) \widehat{\omega}_{j+1}^{(i)}, \quad (3.23)$$

where $g_j(x)$ given by (2.10) denotes the likelihood function as before. Then the μ_{j+1}^N approximated by the indicator measure is given by

$$\mu_{j+1}^N = \sum_{i=1}^N \omega_{j+1}^{(i)} \mathbb{1}_{\mathbb{B}_i}(dx).$$

Note that we choose the mass point $x^{(i)}$ of each box \mathbb{B}_i to calculate $g_j(x)$, i.e., the likelihood function. It is a reasonable choice to approximate the likelihood function of the points in the \mathbb{B}_i . In both prediction step and analysis step, they only evolve weights $\{\omega_j^{(i)}\}_{i=1}^N$ into $\{\omega_{j+1}^{(i)}\}_{i=1}^N$ via $\{\widehat{\omega}_{j+1}^{(i)}\}_{i=1}^N$, and provide a transform from μ_j^N to μ_{j+1}^N . The complete procedure is summarized in Algorithm 2, named Perron-Frobenius operator filter. The algorithm consists of two phases: offline phase to compute P^N by Ulam's method, and online phase to update the approximation of the posterior measure. Besides, the standard Ulam's method becomes inefficient in high-dimensional dynamical systems due to the curse of dimensionality. For this case, we may use the sparse Ulam method [27] instead. It constructs an optimal approximation subspace and costs less computational effort than the standard Ulam's method when a certain accuracy is imposed. Algorithm 1 Perron-Frobenius operator filter

Offline: Compute P^N by Ulam's method Online: 1: Set j = 0 and $\mu_0^N(dx_0) = \mu_0(dx_0)$, compute $\omega_0^{(i)} = \frac{\int_{\mathbb{B}_i} \mu_0 dx_0}{\mu(\mathbb{B}_i)}$ 2: Let $\mathbf{W}_j = [\omega_j^{(1)}, \cdots, \omega_j^{(N)}]$, compute $\widehat{\mathbf{W}}_{j+1} = \mathbf{W}_j P^N$ 3: Define $\widehat{\mu}_{j+1}^N = \sum_{i=1}^N \widehat{\omega}_{j+1}^{(i)} \mathbb{1}_{\mathbb{B}_i}(x)$ 4: Denote $\omega_{j+1}^{(i)}$ by (3.23), define $\mu_{j+1}^N = \sum_{i=1}^N \omega_{j+1}^{(i)} \mathbb{1}_{\mathbb{B}_i}(x)$ 5: $j+1 \rightarrow j$ 6: Go to step 2

3.2 Analysis of error estimate

We analyze the error estimate of the Perron-Frobenius operator filter in this section to explore the factors, which determine convergence of the algorithm. Define a total-variation distance $d(\cdot, \cdot)$ between two probability measures μ and ν as follows:

$$d(\mu, \nu) = \frac{1}{2} \sup_{|f|_{\infty} \le 1} |\mathbb{E}^{\mu}(f) - \mathbb{E}^{\nu}(f)|,$$

where $\mathbb{E}^{\mu}(f) = \int_X f(x)\mu(dx)$ for $f \in L^1(X)$ and $|f|_{\infty} = \sup_x |f(x)|$. The distance $d(\cdot, \cdot)$ can also be characterized by the L^1 norm of the difference between the two PDFs ρ and ρ' , which correspond to the measure μ and ν , respectively, i.e.,

$$d(\mu,\nu) = \frac{1}{2} \int_{X} |\rho(x) - \rho'(x)| dx.$$
(3.24)

The distance induces a metric. To estimate the error, we recall the iteration (3.14) and see that the approximation error of the probability comes from the operator \mathcal{P}^N . To do this, we need the following lemmas.

Lemma 3.1. (Theorem 4.8 in [23]) Suppose that \mathcal{P} is the Perron-Frobenius operator defined in (2.8). Let μ and ν be two arbitrary probability measures. Then

$$d(\mathcal{P}\mu, \mathcal{P}\nu) \le d(\mu, \nu).$$

Lemma 3.2. (Lemma 4.9 in [23]) Let g_j be the likelihood function defined by (2.10) and the operator L_j defined by (2.12). Assume that there exists $\kappa \in (0, 1]$ such that for all $x \in X$ and $j \in \mathbb{N}$,

$$\kappa \le g_j(x) \le \kappa^{-1}.\tag{3.25}$$

Then we have

$$d(L_j\mu, L_j\nu) \le 2\kappa^{-2}d(\mu, \nu)$$

Lemma 3.3. (Theorem 2.4.1 in [28]) Let \mathcal{C}_a be discrete Lipschitz cone defined as

$$\mathcal{C}_a = \{\phi : \frac{\phi(x)}{\phi(y)} \le e^{a|x-y|}, \forall x, y \in \mathcal{R}\}.$$

For each N > 0, the π_N given by (3.16) denotes the projection of L^1 onto V_N . Then for any function $f \in C_a$,

$$||f - \pi_N f||_{L^1} \le (e^{a/N} - 1)||f||_{L^1}.$$

By the above three lemmas, we analyze total-variance distance between the approximate measure μ_J^N and the true measure μ_J and give the following theorem.

Theorem 3.4. If $g_j(x)$ satisfies the condition (3.25) and the probability density ρ_j^N of the measure μ_j^N satisfies $\mathcal{P}\rho_j^N \in \mathcal{C}_a$, $\forall j \in \mathbb{N}$, then

$$d(\mu_J^N, \mu_J) \le \sum_{j=1}^J (2\kappa^{-2})^j \frac{e^a - 1}{2N}.$$

Proof. From the formula (2.13) and (3.14), we apply the triangle inequality to the distance $d(\mu_{i+1}^N, \mu_{j+1})$ and get

$$d(\mu_{j+1}^N, \mu_{j+1}) = d(L_j \mathcal{P}^N \mu_j^N, L_j \mathcal{P} \mu_j)$$

$$\leq d(L_j \mathcal{P}^N \mu_j^N, L_j \mathcal{P} \mu_j^N) + d(L_j \mathcal{P} \mu_j^N, L_j \mathcal{P} \mu_j).$$

According to Lemma 3.2 and Lemma 3.1, it follows that

$$d(\mu_{j+1}^{N}, \mu_{j+1}) \leq 2k^{-2} \left[d(\mathcal{P}^{N} \mu_{j}^{N}, \mathcal{P} \mu_{j}^{N}) + d(\mathcal{P} \mu_{j}^{N}, \mathcal{P} \mu_{j}) \right] \\ \leq 2k^{-2} \left[d(\mathcal{P}^{N} \mu_{j}^{N}, \mathcal{P} \mu_{j}^{N}) + d(\mu_{j}^{N}, \mu_{j}) \right],$$
(3.26)

Let us consider $d(\mathcal{P}^N \mu_j^N, \mathcal{P} \mu_j^N)$. Suppose that ρ'_{j+1} is density function associated with the measure $\mathcal{P}\mu_j^N$. Let ρ_j^N and ρ_{j+1}^N be the density functions of μ_j^N and μ_{j+1}^N , respectively. By the definition of total-variance distance in (3.24), we have

$$d(\mathcal{P}^N \mu_j^N, \mathcal{P} \mu_j^N) = \frac{1}{2} \int_X |\rho_{j+1}^N(x) - \rho_{j+1}'(x)| dx$$

$$= \frac{1}{2} \int_X |\mathcal{P}^N \rho_j^N(x) - \mathcal{P} \rho_j^N(x)| dx$$

$$= \frac{1}{2} \int_X |\pi_N \circ \mathcal{P} \rho_j^N(x) - \mathcal{P} \rho_j^N(x)| dx,$$

where we have used the equation $\mathcal{P}^N = \pi_N \circ \mathcal{P}$ in the last equality. Since $\mathcal{P}\rho_j^N(x) \in \mathcal{C}_a$, we use Lemma 3.3 and get

$$d(\mathcal{P}^{N}\mu_{j}^{N}, \mathcal{P}\mu_{j}^{N}) \leq \frac{1}{2}(e^{a/N} - 1)$$

$$\leq \frac{1}{2N}(e^{a} - 1).$$
(3.27)

With the fact that $\mu_0^N = \mu_0$, we combine (3.27) with (3.26) and repeat the iterating to complete the proof.

Theorem 3.4 estimates the online error of the PFOF algorithm. Since the Perron-Frobenious operator is numerically approximated offline by the matrix form P_n^N given by (3.18), we will analyze the offline error generated by the approximation. Each coefficient of P_n^N is computed by the Monte-Carlo approximation of (3.17) using a set of the sampling points $\{x_i^l\}_{l=1}^n$. We show that the matrix P_n^N converge to the matrix P^N . **Proposition 3.5.** If the matrix P_n^N is defined by (3.18) and P^N is defined by (3.17), then the following convergence in distribution holds,

$$\sqrt{n}((P_n^N)_{ij} - (P^N)_{ij}) \xrightarrow[n \to \infty]{\mathcal{D}} \mathcal{N}(0, \sigma_{ij}^{n,N}), \qquad (3.28)$$

where

$$(\sigma_{ij}^{n,N})^2 = \int_X (\mathbb{1}_{\Psi_\tau^{-1}(\mathbb{B}_j)} \cdot \mathbb{1}_{\mathbb{B}_i})^2 d\mu - (\int_X \mathbb{1}_{\Psi_\tau^{-1}(\mathbb{B}_j)} \cdot \mathbb{1}_{\mathbb{B}_i} d\mu)^2,$$
(3.29)

and $\mathcal{N}(0, \sigma_{ij}^{n,N})$ is the normal distribution with the mean 0 and standard deviation $\sigma_{i,j}^{n,N}$. *Proof.* Note that the entries of P_n^N are given by

$$P_{n,ij}^N = \frac{1}{n} \sum_{l=1}^n \mathbb{1}_{\mathbb{B}_j}(\Psi_\tau(x_i^l)),$$

which is the Monte-Carlo approximation of

$$P_{ij}^N = \frac{\int_X \mathbb{1}_{\Psi_{\tau}^{-1}(\mathbb{B}_j)} \cdot \mathbb{1}_{\mathbb{B}_i} d\mu}{\int_X \mathbb{1}_{\mathbb{B}_i} d\mu}$$

with sampling points x_i^l drawn independently and uniformly from the box \mathbb{B}_i . The denominator $\int_X \mathbb{1}_{\mathbb{B}_i} d\mu$ normalizes the entries P_{ij}^N so that P^N becomes a right stochastic matrix, with each row summing to 1. The convergence result (3.28) follows directly from the convergence of Monte-Carlo integration [29].

Proposition 3.5 indicates that there exits a constant $C^N(\Psi_{\tau}, \alpha)$ determined by the standard deviation $\sigma_{ij}^{n,N}$ with a given confidence rate $\alpha \in [0, 1)$ such that for *m* large enough, the following estimate holds with probability at least α :

$$\| (P_n^N)_{ij} - (P^N)_{ij} \|_{\infty} \le C^N(\Psi_{\tau}, \alpha) n^{-\frac{1}{2}}.$$
(3.30)

The result shows that the convergence of P_n^N to P^N is in $\mathcal{O}(n^{-\frac{1}{2}})$.

3.3 A low-rank Perron-Frobenius operator filter

In PFOF, we note that the number of blocks increases exponentially with respect to dimensions, resulting in the number of basis functions growing rapidly. Therefore, we propose a low-rank approximation, formed by eigenfunctions of the Perron-Frobenius operator, to represent the density. This approach can effectively reduce the number of the required basis functions. Let ρ still be a probability density function of the dynamical system governed by Ψ . Then it can be written as a linear combination of the independent eigenfunctions φ_i of \mathcal{P} . So

$$\rho(x,t) = \sum_{i=1}^{\infty} a_i \varphi_i(x), \quad a_i \in \mathbb{C}.$$

Suppose that λ_i is the eigenvalue corresponding to the eigenfunction φ_i of \mathcal{P} , then

$$\mathcal{P}\rho(x,t) = \sum_{i=1}^{\infty} \lambda_i a_i \varphi_i(x).$$

Actually, the eigenfunction of the discretized PFO can be determined by the following proposition.

Proposition 3.6. Let $B = \{\mathbb{B}_1, \dots, \mathbb{B}_N\} \subset \mathcal{B}$ be a uniform partition of the phase space X. If ξ is the left eigenvector of P^N corresponding to the eigenvalue λ , then λ is also the eigenvalue of the restricted operator $\pi_N \mathcal{P}$ with the eigenfunction $\varphi \triangleq \xi^T \mathbf{U}$, where $\mathbf{U} = [\mathbb{1}_{\mathbb{B}_1}(x), \dots, \mathbb{1}_{\mathbb{B}_N}(x)]^T$.

Proof. Let $\varphi = \sum_{i} \xi^{(i)} \mathbb{1}_{\mathbb{B}_{i}}$. From Eq. (3.21) and Eq. (3.22),

$$\pi_N \mathcal{P} \varphi = \sum_j^N \sum_i^N \xi^{(i)} P_{ij}^N \mathbb{1}_{\mathbb{B}_j}.$$

Since $\xi P^N = \lambda P^N$, i.e., $\lambda \xi^{(j)} = \sum_i \xi^{(i)} P^N_{ij}, \ \forall j \in \mathbb{N}$, we get

$$\pi_N \mathcal{P} \varphi = \sum_j^N \lambda \xi^{(j)} \mathbb{1}_{\mathbb{B}_j} = \lambda \varphi$$

Thus, λ is also an eigenvalue of the restricted operator $\pi_N \mathcal{P}$ with eigenfunction φ .

In order to obtain the spectral expansion of the density function $\rho_j := \rho(x, t_j)$, we define the matrix $\boldsymbol{\varphi} = [\varphi_1, \cdots, \varphi_N]^T$, where $\{\varphi_i\}_{i=1}^N$ are the eigenfunctions with respect to eigenvalues $\{\lambda_i\}_{i=1}^N$, with $|\lambda_1| = 1 \ge |\lambda_2| \ge \cdots \ge |\lambda_N| \ge 0$. Let $\rho_0 = \mathbf{W}_0 \mathbf{U}$ and

$$\Xi = \begin{bmatrix} \xi_1^T \\ \xi_2^T \\ \vdots \\ \xi_N^T \end{bmatrix}.$$

Then the eigenfunction is denoted as $\varphi = \Xi \mathbf{U}$ and the density function of ρ_1 is given by

$$\rho_1 = \pi_N \mathcal{P} \rho_0 = \pi_N \mathcal{P} \mathbf{W}_0 \mathbf{U} = \pi_N \mathcal{P} \mathbf{W}_0 \Xi^{-1} \boldsymbol{\varphi} = \Lambda \mathbf{W}_0 \Xi^{-1} \boldsymbol{\varphi} = \sum_{i=1}^N \lambda_i \varphi_i v_i, \qquad (3.31)$$

where Λ is a diagonal eigenvalue matrix for $\pi_N \mathcal{P}$ and v_i is the column vector of the matrix $V = \mathbf{W}_0 \Xi^{-1}$. The first r major eigenvalues and their corresponding eigenfunctions are used to approximate the density function. If the formula (3.31) is truncated by r < N, then the low-rank model of ρ_1 has the form of

$$\rho_1 = \sum_{i=1}^r \lambda_i \varphi_i v_i.$$

In this way, the low-rank model of density functions at time t_j is given by

$$\rho_j = \sum_{i=1}^r \lambda_i^j \varphi_i v_i, \quad j \in \mathbb{N}.$$

Let us denote the low-rank approximation of Perron-Frobenius operator as $\rho_j = \widetilde{\mathcal{P}}\rho_{j-1} \triangleq \sum_{i=1}^r \lambda_i \varphi_i v_{j-1,i}$, where $v_{j-1,i}$ is the column vector of the matrix $V_{j-1} = \mathbf{W}_{j-1} \Xi^{-1}$. We apply $\widetilde{\mathcal{P}}$ in the Bayesian filter to obtain the low-rank Perron-Frobenius operator filter (lr-PFOF), in which the probability measure satisfies the recursive formula

$$\mu_{j+1}^N = L_j \widetilde{\mathcal{P}} \mu_j^N, \quad \mu_0^N = \mu_0.$$

To describe the following prediction and analysis steps, we first calculate the weak approximation P^N and get the eigenvalues Λ and left eigenvectors Ξ of P^N .

Prediction In this step, we give a model decomposition of the prior density $p(x_{j+1}|Y_j)$. First, the \mathbf{W}_j satisfying $p(x_j|Y_j) = \mathbf{W}_j \mathbf{U}$ is obtained from the previous analysis step. Next, compute the matrix $V_j = \mathbf{W}_j \Xi^{-1}$ and

$$\widehat{\rho}_{j+1} = p(x_{j+1}|Y_j) = \sum_{i=1}^r \lambda_i \varphi_i v_{j,i}.$$

Analysis In this step, we derive the posterior density $p(x_{j+1}|Y_{j+1})$ via Bayes's formula. Multiply $p(x_{j+1}|Y_j)$ by likelihood function g_j and have

$$\rho_{j+1} = p(x_{j+1}|Y_{j+1}) \propto \sum_{i=1}^r \lambda_i \varphi_i v_{j,i} g_j$$

To normalize ρ_{j+1} , we rewrite the $\hat{\rho}_{j+1}$. Since $\varphi_i = \xi_i^T \mathbf{U} = \sum_{k=1}^N \xi_i^{(k)} \mathbb{1}_{\mathbb{B}_k}$, we get

$$\widehat{\rho}_{j+1} = \sum_{i=1}^{r} \lambda_i \sum_{k=1}^{N} \xi_i^{(k)} \mathbb{1}_{\mathbb{B}_k} v_{j,i} = \sum_{k=1}^{N} \left(\sum_{i=1}^{r} \lambda_i \xi_i^{(k)} v_{j,i} \right) \mathbb{1}_{\mathbb{B}_k}$$

Then we multiply by $g_i(x)$ and make a normalization to the weights of $\mathbb{1}_{\mathbb{B}_k}$, such that

$$\omega_{j+1}^{(k)} = \widetilde{\omega}_{j+1}^{(k)} / (\sum_{n=1}^{N} \widetilde{\omega}_{j+1}^{(n)}), \quad \widetilde{\omega}_{j+1}^{(k)} = \sum_{i=1}^{r} \lambda_i \xi_i^{(k)} v_{j,i} g_j(x^{(k)}), \tag{3.32}$$

where $x^{(k)}$ is still the mass point of each box \mathbb{B}_k . The posterior density becomes

$$\rho_{j+1} = \sum_{k=1}^{N} \omega_{j+1}^{(k)} \mathbb{1}_{\mathbb{B}_k} = \mathbf{W}_{j+1} \mathbf{U}.$$

Remark 3.1. Note that the complex eigenvalues and eigenvectors may appear in the eigendecomposition of the matrix P^N . When the stationary distribution $\tilde{\pi}$ of the system satisfies detailed balance, a symmetrization method is designed in [30] to solve the problem. Since Algorithm 2 low-rank Perron-Frobenius operator filter

Offline:

Compute P^N and its eigenvalue Λ and left eigenvector Ξ . Give the eigenfunction $\varphi = \Xi \mathbf{U}$. **Online:**

- 1: Set j = 0 and $\rho_0 = \mathbf{W}_0 \mathbf{U}$, compute $\omega_0^{(i)} = \frac{\int_{\mathbb{B}_i} \mu_0 dx_0}{\mu(\mathbb{B}_i)}$
- 2: Denote $V_j = \mathbf{W}_j \Xi^{-1}$, compute $\widehat{\rho}_{j+1} = \sum_{i=1}^r \lambda_i \varphi_i v_{j,i}$ 3: Define g_j by (2.10), give $\rho_{j+1} \propto \sum_{i=1}^r \lambda_i \varphi_i v_{j,i} g_j$
- 4: Normalize weights by (3.32) and obtain \mathbf{W}_{j+1} , let $\rho_{j+1} = \sum_{k=1}^{N} \omega_{j+1}^{(k)} \mathbb{1}_{\mathbb{B}_k}$.
- 5: $j+1 \rightarrow j$
- 6: Go to step 2

 ρ_0, ρ_1, \cdots can be seen as a Markov chain with transition matrix P^N , we suppose that P^N satisfies detailed balance with respect to $\tilde{\pi}$, i.e.,

$$\widetilde{\pi}_i P_{ij}^N = \widetilde{\pi}_j P_{ji}^N, \quad \forall i, j \in N.$$

Then P^N can be symmetrized by a similarity transformation

$$S = \widetilde{\Lambda} P^N \widetilde{\Lambda}^{-1}, \quad \text{where } \widetilde{\Lambda} = \begin{bmatrix} \sqrt{\widetilde{\pi}_1} & & \\ & \sqrt{\widetilde{\pi}_2} & \\ & & \ddots & \\ & & & \sqrt{\widetilde{\pi}_N} \end{bmatrix}$$

Here the S is a symmetric matrix and this can be easily checked by detailed balance equation. It is known that S has a full set of real eigenvalues $\alpha_j \in \mathbb{R}$ and an orthogonal set of eigenvectors w_j . Therefore, P^N has the same eigenvalues α_j and real left eigenvectors

$$\psi_j = \Lambda w_j$$

3.4 Extension to continuous-time filtering problems

In this subsection, we consider a continuous-time filtering problem, where the state model and observation are by the following SDEs,

$$\frac{dx}{dt} = f(x) + \sqrt{\Sigma_c} \frac{dW_t}{dt}, \quad x(t_0) \sim \mathcal{N}(m_0, C_0), \tag{3.33}$$

$$\frac{dz}{dt} = h(x) + \sqrt{R_c} \frac{dW_t}{dt}, \quad z(0) = 0.$$
(3.34)

Here Σ_c is the covariance of model error and R_c is the covariance of observation error. Suppose that the posterior measure μ_t governed by the continuous-time problem has Lebesgue density $\rho(\cdot,t): \mathbb{R}^n \to \mathbb{R}^+$ for a fixed t. Let $\rho(x,t) = r(x,t) / \int_{\mathbb{R}^n} r(x,t) dx$, where r is the unnormalized density. For a positive definite symmetric matrix $A \in \mathbb{R}^{p \times p}$, we define the weighted inner product $\langle \cdot, \cdot \rangle_A = \langle A^{-\frac{1}{2}}, A^{-\frac{1}{2}} \rangle$ on the space $L^2([0,T];\mathbb{R}^p)$. The resulting norm $|\cdot|_A = |A^{-\frac{1}{2}} \cdot|$. In the continuous filtering problem, our interest is to find the distribution of the random variable $x(t)|\{z(s)\}_{s\in[0,t]}$ as the time t increases. Zakai stochastic partial differential equation (SPDE) is a well-known equation whose solution characterizes the unnormalized density of posterior distribution [35]. The Zaikai equation has the form of

$$\frac{\partial r}{\partial t} = \mathscr{A}_{PF}r + r\left\langle h, \frac{dz}{dt} \right\rangle_{R_c}.$$
(3.35)

The partial differential operator \mathscr{A}_{PF} generates a continuous Perron-Frobenius semigroup $\{\mathcal{P}_t, t \geq 0\}$. Let $\{\mathcal{Q}_s^t, 0 \leq s \leq t\}$ be the stochastic semigroup [31] associated with with the following SDE

$$\frac{dr'}{dt} = r' \left\langle h, \frac{dz}{dt} \right\rangle_{R_c}.$$
(3.36)

Then the Zakai equation (3.35) can be approximated by the following Trotter-like product formula

$$r_{j+1} = Q_{t_j}^{t_{j+1}} \mathcal{P}_\tau r_j, \qquad (3.37)$$

where $\tau = t_{j+1} - t_j$, $\forall j \in \mathbb{N}$. For the fixed τ , \mathcal{P}_{τ} is still denoted by \mathcal{P} . By the reference [31], the $Q_{t_j}^{t_{j+1}}$ describes the solution of the equation (3.36), i.e.,

$$Q_{t_j}^{t_{j+1}}r(x) = \exp\left(\langle h(x), z_{j+1} - z_j \rangle_{R_c} - \frac{\tau}{2} |h(x)|_{R_c}^2\right) r(x).$$

With the discrete scheme (3.37), we utilize the Perron-Frobenius operator to solve Zakai equation, rather than using Fokker-Planck operator \mathscr{A}_{PF} . Thus, we discretize \mathcal{P} by Ulam's method and project the density function onto V_N . Let P^N be the discretization of \mathcal{P} . Let \mathbf{W}_j and \mathbf{W}_{j+1} be the weights vectors with respect to $\pi_N r_j$ and $\pi_N r_{j+1}$. Denote $g_j^c(x) = \exp(\langle h(x), z_{j+1} - z_j \rangle_{R_c} - \frac{\tau}{2} |h(x)|_{R_c}^2)$ and

$$\mathbf{G}_{j} = \begin{bmatrix} g_{j}^{c}(x^{(1)}) \\ g_{j}^{c}(x^{(2)}) \\ \vdots \\ g_{j}^{c}(x^{(N)}) \end{bmatrix},$$

where $x^{(i)}$ is the mass point of \mathbb{B}_i . Then the transition of density functions turns into a map of the weights,

$$\mathbf{W}_{j+1} = \mathbf{G}_j \odot \left(\mathbf{W}_j P^N \right)$$

Here \odot denotes Hadamard product. In this case, the PFO is extended to the continuous-time filtering problem to estimate the posterior density function.

4 Comparison with particle filter

Particle filter (PF) [32, 33] is an important filtering method to sequentially approximate the true posterior filtering distribution $p(x_i|Y_i)$ in the limit of a large number of particles. In

practice, we approximate the probability density by a combination of locations of particles and weights associated with Dirac functions. Particle filter proceeds by varying the weights and determining the particle Dirac measures. It is able to take care of non-Gaussian and nonlinear models. In this section, we will compare the computational accuracy and differences between PFOF and PF.

Accordingly, we define μ_j^N as the posterior empirical measure on \mathbb{R}^N approximating truth posterior probability measure μ_j and define $\hat{\mu}_j^N$ on \mathbb{R}^N as the approximation of the prior probability measure $\hat{\mu}_j$. Let

$$\mu_j \approx \mu_j^N := \sum_{n=1}^N \omega_j^{(n)} \delta_{x_j^{(n)}}, \quad \widehat{\mu}_{j+1} \approx \widehat{\mu}_{j+1}^N := \sum_{n=1}^N \widehat{\omega}_{j+1}^{(n)} \delta_{\widehat{x}_{j+1}^{(n)}},$$

where $x_j^{(n)}$ and $\hat{x}_{j+1}^{(n)}$ are particle positions, and $\omega_j^{(n)} > 0$, $\hat{\omega}_{j+1}^{(n)} > 0$ are the associated weights satisfying $\sum_{n=1}^{N} \omega_j^{(n)} = 1$, $\sum_{n=1}^{N} \hat{\omega}_{j+1}^{(n)} = 1$. The empirical distribution is completely determined by particle positions and weights. The objective of particle filter is to calculate the update $\{x_j^{(n)}, \omega_j^{(n)}\} \rightarrow \{\hat{x}_{j+1}^{(n)}, \hat{\omega}_{j+1}^{(n)}\}$ and $\{\hat{x}_{j+1}^{(n)}, \hat{\omega}_{j+1}^{(n)}\} \rightarrow \{x_{j+1}^{(n)}, \omega_{j+1}^{(n)}\}$, which define the prediction step and analysis step, respectively. Monte-Carlo sampling is used to determine particle positions in the prediction and Bayesian rule is used to update of the weights in the analysis.

Prediction In this step, the prediction phase is approximated by the Markov chain $\{\Psi(x_j)\}_{j\in\mathbb{N}}$ with transition kernel $p(x_j, x_{j+1}) = p(x_{j+1}|x_j)$. We draw $\widehat{x}_{j+1}^{(n)}$ from the kernel p started from $x_j^{(n)}$, i.e., $\widehat{x}_{j+1}^{(n)} \sim p(x_j^{(n)}, \cdot)$. We keep the weights unchanged so that $\widehat{\omega}_{j+1}^{(n)} = \omega_j^{(n)}$, and obtain the prior probability measure

$$\widehat{\mu}_{j+1}^N = \sum_{n=1}^N \omega_j^{(n)} \delta_{\widehat{x}_{j+1}^{(n)}}$$

Analysis In this step, we apply Bayes's formula to approximate the posterior probability measure. To do this, we fix the position of the particles and update the weights. With the definition of $g_i(x)$ in (2.10), we have the empirical posterior distribution

$$\mu_{j+1}^N = \sum_{n=1}^N \omega_{j+1}^{(n)} \delta_{\widehat{x}_{j+1}^{(n)}}$$

where

$$\omega_{j+1}^{(n)} = \widetilde{\omega}_{j+1}^{(n)} / (\sum_{n=1}^{N} \widetilde{\omega}_{j+1}^{(n)}), \quad \widetilde{\omega}_{j+1}^{(n)} = g_j(\widehat{x}_{j+1}^{(n)}) \omega_j^n.$$
(4.38)

The first equation in (4.38) is a normalization. Sequential Importance Resampling (SIR) particle filter is a basic particle filter and shown in Algorithm 1. A resampling step is introduced in the algorithm. In this way, we can deal with the initial measure μ_0 when it is not a combination of Dirac functions. We can also deal with the case when some of the particle weights are close to 1. The algorithm shows that each particle moves according

to the underlying model and is reweighted according to the likelihood. By the iteration of Bayesian filtering , we rewrite the particle filter approximated by the form

$$\mu_{j+1}^N = L_j S^N \mathcal{P} \mu_j^N, \quad \mu_0^N = \mu_0, \tag{4.39}$$

where the operator S^N is defined as follows:

$$(S^N \mu)(dx) = \frac{1}{N} \sum_{n=1}^N \delta_{x^{(n)}}(dx), \quad x^{(n)} \sim \mu \quad \text{i.i.d.}.$$

Algorithm 3 Sequential Importance Resampling particle filter

1: Set j = 0 and $\mu_0^N(dx_0) = \mu_0(dx_0)$ 2: Draw $x_j^{(n)} \sim \mu_j^N, n = 1, \dots, N$ 3: Set $\omega_j^{(n)} = 1/N, n = 1, \dots, N$, redefine $\mu_j^N := \sum_{n=1}^N \omega_j^{(n)} \delta_{x_j^{(n)}}$ 4: Draw $\widehat{x}_{j+1}^{(n)} \sim p(x_j^{(n)}, \cdot)$ 5: Define $\omega_{j+1}^{(n)}$ by (3.23) and $\mu_{j+1}^N = \sum_{n=1}^N \omega_{j+1}^{(n)} \delta_{\widehat{x}_{j+1}^{(n)}}$ 6: $j+1 \rightarrow j$ 7: Go to step 2

By (4.39), we find that the randomness for the probability measure is caused by the sampling operator S^N and the convergence of particle filter depends on the number of particles. The particle filter does recover the truth posterior distribution as the number of particles tends to infinity [34]. The following theorem gives a convergence result for PF.

Theorem 4.1. (Theorem 4.5 in [23]) Let m be the number of particles and μ_j^m the approximation measure in SIR particle filter. Assume that $\kappa \in (0, 1]$ is the constant defined in Lemma 3.2, then the total-variance distance between μ_J^m and μ_J is estimated by

$$d(\mu_J^m, \mu_J) \le \sum_{j=1}^J (2\kappa^{-2})^j \frac{1}{\sqrt{m}}.$$
(4.40)

Let J be fixed in Theorem 3.4 and Theorem 4.1. We find that the convergence rate of particle filter depends on the number of particles m. Similarly, the convergence of PFOF is determined by the number of blocks N used in the Ulam's method. When N = m, i.e., the same number of basis functions in the two methods, the rate of convergence is $\mathcal{O}(\frac{1}{N})$ in PFOF and $\mathcal{O}(\frac{1}{\sqrt{m}})$ in SIR particle filter. The analysis shows that PFOF converges faster than the particle filter.

Sampling from high-dimensional and complex transition kernels is difficult to realize in PF. The PFOF avoids the sampling and uses a data-driven approximation instead, which requires short-term path simulations rather than the form of transition density. Particle degeneracy is also a significant issue. As the number of effective particles decreases gradually, the efficiency of the particle filter becomes worse.

It is known that particle filter is inefficient for high-dimensional models because of degeneracy. So the accurate estimate of posterior PDF requires a great number of particles that scales exponentially with the size of the system. In addition to resampling, adding jitter and localisation are effective modifications to solve the problem. The PFOF also has the "curse of dimensionality" problem in high dimensions as the partition scale expansion. One solution to circumvent this problem is the sparse Ulam method. The low-rank Perron-Frobenius operator filter can enhance the efficiency of filtering problems.

5 Numerical results

In this section, we present some numerical examples for filtering problems using the proposed PFOF. The system dynamics is unknown and some observations are given in the filtering problems. The PFOF and lr-PFOF are implemented to estimate posterior PDFs of the stochastic filtering problems. In Section 5.1, we consider an Ornstein-Uhlenbeck (O-U) process to identify the Gaussian PDF of the system and estimate its posterior PDFs with observations known. In Section 5.2, we consider a nonlinear filtering problem governed by Beneš SDE, and estimate the non-Gaussian posterior PDFs. In Section 5.3, we consider a continuous-time filtering problem, which is a classical chaotic system Lorenz'63 model with observations, to model posterior density of the state. We compare the proposed PFOF/lr-PFOF with particle filter and Extended Kalmn filter (ExKF). Numerical results show that PFOF achieves a better posterior PDF estimates than PF, and a more accurate state estimates than ExKF.

5.1 O-U process

Let us consider an O-U process, which is a one-dimensional linear dynamical system,

$$dx_t = -\lambda x_t dt + dW_t, \quad x(t_0) \sim \mathcal{N}(m_0, C_0),$$

where $\lambda > 0$ and W_t is a standard Brownian motion. We now consider a state-space model formed by a discretization of the O-U process and the discrete observations of the state as follows,

$$\begin{cases} x(t_{k+1}) = \exp(-\lambda \Delta t_k) x(t_k) + q_k, \quad q_k \sim \mathcal{N}(0, \Sigma_k), \\ y(t_k) = H x(t_k) + r_k, \quad r_k \sim \mathcal{N}(0, R), \end{cases}$$
(5.41)

where $\Sigma_k = \exp(-2\lambda\Delta t_k)$, H = I and $R = \sigma^2$. The parameters are given by $\lambda = 1/2$, $m_0 = 2, C_0 = 0.1$ and $\sigma = 1$. To apply PFOF, we compute Perron-Frobenius operator P_{τ} using Ulam's method with time step $\tau = 0.1$ and obtain an approximation form $P_{\tau}^N \in \mathbb{R}^{N \times N}$ of \mathcal{P}_{τ} . We take the phase space of x_t is [-6, 6] and divide it into N = 100 grids, and each interval $[z_k, z_{k+1}], k = 0, \dots, N - 1$, defines a box \mathbb{B}_k . We define an indicator function $\mathbb{1}_{\mathbb{B}_k}(x)$ on each \mathbb{B}_k and randomly choose n = 100 sample points in the box to calculate P_{τ}^N . Given initial Gaussian distribution $\mathcal{N}(2, 0.1)$, we rewrite μ_0 as a vector W_0 , which denotes the coefficients of μ_0^N . The P_{τ}^N acts on the weight vector to estimate probability value of x_t on each \mathbb{B}_k , i.e., $\mathbb{P}(x_t \in \mathbb{B}_k), t = q\tau, q = 0, 1, 2 \cdots$. Thus, we get the discrete probability density function (PDF) of x_t at t. The simulation PDFs at different times are shown in the left column of Figure 5.1. By the figure, we see that the PDFs estimated by PFO are close



Figure 5.1: The prior PDF estimated by PFO (left column), posterior PDF by PFOF (middle column) and posterior PDF by particle filter (right column) at different times.

to the truth. By this way, the PDF is computed without solving Fokker-Planck equation and the estimation of PDF is actually the prior density in the model (5.41).

Then we compute posterior probability density of the state-space model (5.41). We set N = 500 and n = 100. The posterior probability density function is estimated by Algorithm 1 and the results are displayed in the middle column of Figure 5.1. From Figure 5.1, we find that the empirical posterior PDFs estimated by PFOF are close to the Gaussian posterior densities. To make comparison with PFOF, the particle filter is also used for the filtering problem. In the particle filter, 500 particles are drawn randomly to generate Dirac measure and construct empirical measure. Thus, the number of basis functions is equal to each other in the two methods. Figure 5.1 clearly shows that the empirical PDF calculated by PFOF is more accurate than that by PF. The numerical results support Theorem 3.4 and Theorem 4.1.

5.2 Beneš-Daum filter

In this subsection, we apply PFOF to a nonlinear filtering problem, whose state-space model is defined by the Beneš stochastic difference equation,

$$dx_t = \tanh(x_t)dt + dW_t, \tag{5.42}$$

with initial condition $x_0 = 0$. Refer to [36], the probability density function of the equation (5.42) is given by

$$p(x(t)) = \frac{1}{\sqrt{2\pi t}} \frac{\cosh(x(t))}{\cosh(x_0)} \exp\left(-\frac{t}{2}\right) \exp\left(-\frac{1}{2t}(x(t) - x_0)\right).$$

We take the phase space [-15, 15] and uniformly divide it into 100 (N = 100) grids $[z_k, z_{k+1}]$, k = 0, ..., N - 1, each of which corresponds to a box \mathbb{B}_k . The Ulam's method is used to approximate PFO. The time step is set as $\tau = 0.5$ and the number of random sample points m = 400. The predicted PDF of x_t at t = 1, t = 2.5 and t = 5 are shown in Figure 5.2. The PDFs are separately estimated by discretized PFO matrix $P^N \in \mathbb{R}^{100 \times 100}$ and low-rank approximation of PFO with truncation r = 30. We see that PDFs at t = 2.5 and t = 5 have two modes and the PFO can fairly approximate the two modes.



Figure 5.2: The PDF estimated by PFO and low-rank model at different times.

First we want to calculate the truth posterior filtering distribution of the model (5.42) subject to observation. In this example, the observation model satisfies

$$p(y_k|x(t_k)) = \mathcal{N}(y_k|x(t_k), \sigma^2).$$
(5.43)

According to [36] (Chapter 10.5), the transition density of the Beneš SDE is given by

$$p(x(t_k)|x(t_{k-1})) = \frac{1}{\sqrt{2\pi\Delta t_{k-1}}} \frac{\cosh(x(t_k))}{\cosh(x(t_{k-1}))} \exp\left(-\frac{1}{2}\Delta t_{k-1}\right) \times \exp\left(-\frac{1}{2\Delta t_{k-1}}(x(t_k) - x(t_{k-1}))^2\right),$$

where $\Delta t_{k-1} = t_k - t_{k-1}$. If we assume that the filtering solution at time t_{k-1} is of the form

$$p(x(t_{k-1})|y_{1:k-1}) \propto \cosh(x(t_{k-1})) \exp\left(-\frac{1}{2P_{k-1}}(x(t_{k-1})-m_{k-1})^2\right)$$

for given m_{k-1} and P_{k-1} . Then we use the Chapman-Kolmogorov equation and give the prior density

$$p(x(t_k)|y_{1:k-1}) \propto \cosh(x(t_k)) \exp\left(-\frac{1}{2P_k^-}(x(t_k)-m_k^-)^2\right),$$

where

$$m_k^- = m_{k-1},$$

 $P_k^- = P_{k-1} + \Delta t_{k-1}.$

The m_k^- and P_k^- are sufficient statistics representing prior density functions. By Bayes' formula, the posterior density of $x(t_k)$ is given by

$$p(x(t_k)|y_{1:k}) \propto \cosh(x(t_k)) \exp\left(-\frac{1}{2P_k}(x(t_k)-m_k)^2\right), \qquad (5.44)$$

where the equations of parameters m_k and P_k in the posterior density satisfy

$$m_{k} = m_{k}^{-} + \left(\frac{P_{k}^{-}}{P_{k}^{-} + \sigma^{2}}\right)(y_{k} - m_{k}^{-}),$$
$$P_{k}^{-} = P_{k-1} + \Delta t_{k-1}.$$

Thus, the reference posterior distribution is defined by (5.44).

To apply PFOF to the nonlinear filtering problem, we make a finer division of the phase interval [-15, 15] to obtain 400 boxes. Besides, we choose enough sample points in Ulam's method to reduce error of Monte-Carlo as much as possible. The observations y_k are artificially obtained by simulating the underlying model (5.42) and adding noise according to (5.43), where $\sigma = 1$. The observable interval is [0,5] with a time step $\Delta t_k = 0.1$. The initial distribution for the filtering process is chosen to be $m_0 = 0$, $P_0 = 2$. Particularly, we also use the particle filter as a comparison. In the prediction, we are not allowed to draw sample points directly because of a complex transition probability density function. We use Acceptance-Rejection method to resolve the issue. We first show the results of posterior mean estimated by PFOF and lr-PFOF (r=40) in Figure 5.3, together with truth and observations. The mean is obtained by averaging the posterior distribution of PFOF/lr-PFOF and it is close to the truth as the figure shows.

The posterior densities estimated by PFOF, lr-PFOF and particle filter are shown in Figure 5.4 together with the truth. The truncation parameters in lr-PFOF are separately set as r = 10, r = 20 and r = 40. The estimation accuracy of lr-PFOF gradually improves as the number of truncation basis functions increases, and achieves almost the same as PFOF when r = 40 < N = 400. Although the number of basis functions is the same in both PFOF and particle filter, there exit clear difference between the two methods. The results show that the accuracy of PFOF is higher than that of particle filter in the non-Gaussian and nonlinear filtering problem. This further confirms the theoretical analysis in Section 3. As shown in Table 1, both PFOF and lr-PFOF use less CPU-time than SIR particle filter does. Actually, the CPU-time in particle filter is mainly from Acceptance-Rejection sampling. From the table, it can be seen that lr-PFOF can reduce online computation time comparing to PFOF.



Figure 5.3: The mean estimated by PFOF and lr-PFOF.



Figure 5.4: The posterior PDF by lr-PFOF (left column), PFOF (middle column) and particle filter (right column) at different times.

Table 1: CPU-time (seconds) for posterior PDF with different methods.

Methods	PFOF	lr-PFOF $(r=10)$	lr-PFOF (r= 20)	lr-PFOF (r=40)	particle filter
offline	0.1599	0.2536	0.2649	0.2689	- 6906.9486
online	0.0673	0.0150	0.0431	0.0641	

5.3 Lorenz'63 model

Lorenz developed a mathematical model for atmospheric convection in 1963. The Lorenz'63 model is the simplest continuous-time system to exhibit sensitivity to initial conditions and chaos, and it is popular example used for data assimilation. For some parameters and initial conditions, the system may perform a chaotic behaviour. The model consists of three coupled nonlinear ordinary differential equations with the solution $v = (v_1, v_2, v_3) \in \mathbb{R}^3$. We consider the Lorenz'63 model with additive white noise,

$$\begin{cases} \frac{dv_1}{dt} = a(v_2 - v_1) + \sigma_1 \frac{dW_1}{dt} \\ \frac{dv_2}{dt} = -av_1 - v_2 - v_1v_3 + \sigma_2 \frac{dW_2}{dt} \\ \frac{dv_3}{dt} = v_1v_2 - bv_3 - b(r+a) + \sigma_3 \frac{dW_3}{dt} \\ v(0) \sim \mathcal{N}(m_0, C_0), \end{cases}$$

where W_j are Brownian motions assumed to be independent. We use the classical parameter values $(a, b, r) = (10, \frac{8}{3}, 28)$ and set $\sigma_1 = \sigma_2 = \sigma_3 = 2$. The initial mean m_0 is given by (0, 0, 0) and covariance matrix is an identity matrix $I_3 \in \mathbb{R}^{3\times 3}$. We give the continuous observation z(t), which is governed by a SDE

$$\begin{cases} \frac{dz}{dt} = h(v) + \gamma \frac{dW_z}{dt}\\ z(0) = 0, \end{cases}$$

with $\gamma = 0.2$. The purpose of this example is to explore the performance of PFOF in continuous-time filtering problems. We compare the assimilation results based on Perron-Frobenius operator and continuous-time Extended Kalman filter. The posterior means estimated by the two methods are shown in Figure 5.5 and Figure 5.7. The two figures are corresponding to different observations h(v) = Hv, where the former is determined by H = [0, 1, 0] and the latter is determined by H = [0, 0, 1]. In particular, we find that the choice of observations in Lorenz models is quite influential, especially for ExKF. The stability of ExKF significantly depends on the observation. Because the insufficient observations may keep the filter away from the truth and cause significant model error, and it may easily lead to the numerical instability once the deviation occurs. However, the results of reconstruction by PFOF much less affected by observation model, so the method shows much better robustness than ExKF.

Figure 5.6 shows the consequence of mean-square error with v_2 or v_3 as the different observation. For ExKF, we find that there is a large error in estimating mean by ExKF when the third component v_3 is observed. To better visualize the results, we compare the trajectories of mean obtained by PFOF and ExKF in Figure 5.8 together with truth. We find the trajectory mean of PFOF agrees with the truth more than the the trajectory mean of ExKF.

For v_3 as an observation, the one-dimensional and two-dimensional marginal probability distributions are displayed in Figure 5.9. The figure aims to intuitively describe distribution of the single value and correlation of the different components. As shown in the figure,



Figure 5.5: The posterior mean of each component by ExKF and PFOF in Lorenz'63 model with continuous observation. The component v_2 is observed.



Figure 5.6: The mean-square error $||v(t) - m(t)||_2^2$ of filters.



Figure 5.7: The posterior mean of each component by ExKF and PFOF in Lorenz'63 model with continuous observation. The component v_3 is observed.



Figure 5.8: The trajectories of mean by PFOF and ExKF.

one-dimensional marginal distributions of the observed component are closer to Gaussian distributions than the other two components. This phenomenon reflects that when a component is used as an observation, its mean estimates will be more accurate than the other unobserved components.

The results above show that PFOF has a higher accuracy for state estimates than ExKF in this chaotic nonlinear system. The former can also give estimates of probability density functions to gain more information of the state in the probabilistic sense.

6 Conclusions

A new filtering method was proposed to estimate filtering distribution of the state under the framework of Perron-Frobenius operator. We formulated filtering problems for discrete and continuous stochastic dynamical systems and applied the Perron-Frobenius operator to propagation of the posterior probability density function. The finite-dimensional approximation of the PFO was realized by Ulam's method, which provides a Galerkin projection space spanned by indicator functions. With Ulam's method, the posterior PDF was discretized and expressed by the weights of basis functions. Then the evolution of PDF became the transition of the weights vectors, which were iterated by PFO and likelihood function. This procedure was called Perron Frobenius operator filter. Thus, the empirical PDF was determined by a convex combination of indicator functions. We gave an error estimate of the proposed method and proved that its accuracy is higher than that of particle filters. Furthermore, a low-rank Perron-Frobenius operator filter was presented to approximate density functions via spectral decomposition. The decomposition was realized by eigendecomposition of discretized PFO. Finally, the proposed method was implemented for three stochastic filtering problems, which included a linear discrete system, a nonlinear discrete system and a nonlinear continuous chaotic system. The numerical results showed that the proposed method has better accuracy and better robustness compared with particle filters and ExKF.

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Figure 5.9: 1-D and 2-D posterior marginal probability density functions of v .

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