
A Variational Approximation for Analyzing the Dynamics of Panel Data

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Abstract

Panel data involving longitudinal measurements of the same set of participants taken over multiple time points is common in studies to understand childhood development and disease modeling. Deep hybrid models that marry the predictive power of neural networks with physical simulators such as differential equations, are starting to drive advances in such applications. The task of modeling not just the observations but the hidden dynamics that are captured by the measurements poses interesting statistical/computational questions. We propose a probabilistic model called ME-NODE to incorporate (fixed + random) mixed effects for analyzing such panel data. We show that our model can be derived using smooth approximations of SDEs provided by the Wong-Zakai theorem. We then derive Evidence Based Lower Bounds for ME-NODE, and develop (efficient) training algorithms using MC based sampling methods and numerical ODE solvers. We demonstrate ME-NODE’s utility on tasks spanning the spectrum from simulations and toy data to real longitudinal 3D imaging data from an Alzheimer’s disease (AD) study, and study its performance in terms of accuracy of reconstruction for interpolation, uncertainty estimates and personalized prediction.

1 INTRODUCTION

Observational studies in the social and health sciences often involve acquiring repeated measurements **over time** for participants/subjects. If most participants stay enrolled, we can consider each row in the *panel* to correspond to longitudinal measurements or records of an individual, at regularly or irregularly sampled time points. Modeling development or

growth trends while accounting for *variability within and across individuals* leads to the need for statistical models for analysis of such “panel data” [Kreindler and Lumsden, 2006, Katsev and L’Heureux, 2003].

The modeling of temporal processes can be set up as a regression task, where a function (with unknown parameters) that is plausible for the domain is estimated using the observed longitudinal data samples. Apart from splines and tools from functional data analysis, a common alternative is to use differential equations [Chen and Wu, 2008, Liang and Yu, 2013, Fang et al., 2011], which provides expressive power and many computational tools developed over decades. However, differential equations do not directly account for variability within and across subjects – hallmark features of panel data. To capture some of these characteristics, the widely used Auto-Regressive Models (ARMA) literature incorporates white noise type functions in differential equation models, leading to various forms of stochastic differential equation (SDE) [Brockwell, 2001, MaCurdy, 1982, Hedeker and Gibbons, 2006],

$$z_t = f_\mu(z, t)dt + L_\Sigma(z, t) \circ d\beta(t) \quad (1)$$

where $z_t \in \mathbf{R}$, f , L denote the drift and noise sensitivity functions with unknown parameters μ and Σ respectively.

Suppose we are given a set of (partial) measurements $\{z_{t_k}\}$ at certain K time points $\{t_k\}_{k=1}^K$, and an efficient numerical scheme to simulate the SDE in (1). Then, the unknown parameters μ and Σ can be found by simply maximizing the likelihood function $p(z_t|\mu; \Sigma)$. In principle, it is straightforward to extend the model in (1) to high dimensional z_t . However, there are two main technical challenges in this setting with stand-alone likelihood based methods: **(i)** such an approach requires a large number of longitudinal measurements which is often infeasible, especially in the applications that motivate our work [Marinescu et al., 2018]; **(ii)** numerical schemes to simulate nonlinear SDEs in high dimensions are quite involved. Indeed, these issues become pronounced when we assume that the observed data is not z but x which are actually measurements governed by a

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process that reflects the dynamics z .

The literature provides a principled way, called **Bayesian filtering**, to tackle the problem described above, see [Särkkä and Solin, 2014], Ch 7. Let us consider that the object or measurement x is evolving as non-linear function D from a *latent* measure of dynamics/progression z_t [Pier-son et al., 2019, Hyun et al., 2016, Whitaker et al., 2017]: $x_t = D(z_t) + \varepsilon_t$. So, it is natural to think of the observable x of an unknown dynamics z – which we can call the “latent” representation. The goal in Bayesian filtering, which aligns nicely with our task, is to compute, $p(z|x_{t_1}, \dots, x_{t_K})$. Interestingly, under some assumptions, closed form solutions for the posterior distributions $p(z|x_{t_1}, \dots, x_{t_K})$ are available, see Chapter 10 in [Särkkä and Solin, 2019]. However, these assumptions are hard to verify in general, and direct utility of these approaches for modern applications is not obvious.

Main ideas/contributions. The most important takeaway from the description above is not the mechanics of *how* Bayesian filtering is carried out in practice, rather, *what* it seeks to estimate. If we focus on the key object of interest – the conditional distribution – we realize that recent works in machine learning do provide a recipe that exploits the universal approximation properties of neural networks to represent fairly complex conditional distributions. In this case, the parameters are simply trained using off-the-shelf procedures, and DE numerical solvers are required for (1). While both ODE and SDE solvers are available, SDE is typically less efficient [Li et al., 2020, Liu et al., 2019]. Notice that when $L_\Sigma \equiv 0$, that is, the observable x follows an ODE and corresponding solvers can be applied, the approach that would instantiate this idea has already been successfully tried in [Yildiz et al., 2019]. However, given the Bayesian filtering motivation above, is there a way to utilize ODE solvers, while preserving stochastic nature of (1)?

Our development begins by rewriting the noise term in (1) with a series of basis functions with standard normal coefficients. We show that this modification enables incorporating random effects in our predictions – which appropriately models the variability of the data, the requirement for successful analysis of panel data. More importantly, using our approach, we show that for a special class of latent SDE based models, which we will refer as Mixed Effect Neural ODE (ME-NODE), the parameters of the underlying neural network can be trained efficiently *without* backpropagating through any SDE solvers. To achieve this, we derive the Evidence Lower Bound loss – where widely available libraries for numerical ODE solvers – are sufficient and directly applicable. We show applications to brain imaging where our formulation can provide personalized prediction together with uncertainty, a feature of Bayesian methods.

2 BACKGROUND/NOTATION

In this section we present the notations and some basic concepts we use in the paper.

Notation. For a time-varying vector $\mathbf{z} = (z_0, \dots, z_n)$, (with n time points) we denote a vector without the j -th component as a $\mathbf{z}_{-j} = (z_0, \dots, z_{j-1}, z_{j+1}, \dots, z_n)$. Often each time point z_t is vector valued, and in the rest of the paper we denote it by $\mathbf{z}_t \in \mathbf{R}^p$, where p is the number of variables. Thus, to denote a time-varying vector where each time point represents a p dimensional vector we use $\mathbf{z} = (\mathbf{z}_0, \dots, \mathbf{z}_n)$, where each $\{\mathbf{z}_t\}_{t=1}^n \subset \mathbf{R}^p$. We denote the indicator function as $\mathbb{1}_{\mathbf{x}}\{\mathbf{y}\}$: it is 1 if $\mathbf{y} = \mathbf{x}$, and 0 otherwise. We refer to a general form of ODE as $\dot{\mathbf{z}}_t = h(\mathbf{z}_t)$, where $h(\mathbf{z}_t)$ defines a “trajectory” and depends on the current value of the process at time t . Without loss of generality, in this section we assume that the DE is defined on \mathbf{R} .

Smooth Approximations of SDE. Consider the standard form of the Stratonovich SDE given by (1). By the Wong and Zakai theorem [Hairer and Pardoux, 2015], the solution to (1) can be approximated asymptotically ($N \rightarrow \infty$) by the solution of the following equation:

$$\dot{z}_t = f(z, t) + L(z, t) \sum_{n=1}^N b_n \varphi_n(t), \quad (2)$$

where $b_n \stackrel{i.i.d.}{\sim} \mathcal{N}(0, 1)$ and $\{\varphi_n\}$ are a suitable set of basis functions. Based on mild simplifying assumptions, we get

$$\dot{z}_t = f(z, t) + g(z, t)b \quad (3)$$

is an approximation of Stratonovich’s SDE in Equation (1), where $b \sim \mathcal{N}(0, 1)$.

Proof. Assume that $\exists g(z, t) < \infty$ and $\forall n, \exists \sigma_n(z, t) < \infty$ such that

$$\varphi_n(t)L(z, t) = \frac{\sigma_n(z, t)}{\sqrt{\sum_{n=1}^N \sigma_n^2(z, t)}} g(z, t). \quad \text{Then,} \quad (4)$$

$$\sum_{n=1}^N \varphi_n(t)L(x, t) = g(z, t) \frac{1}{\sqrt{\sum_{n=1}^N \sigma_n^2}} \sum_{n=1}^N \xi_n(z, t),$$

where $\xi_n(z, t) = b_n \sigma_n(z, t) \sim \mathcal{N}(0, \sigma_n(z, t))$. If for some $\delta > 0$,

$$\frac{1}{\left(\sqrt{\sum_{n=1}^N \sigma_n^2}\right)^{2+\delta}} \sum_{n=1}^N \mathbb{E} \left(|\xi_n - \mu_n|^{2+\delta} \right) \xrightarrow{N \rightarrow \infty} 0, \quad (5)$$

then according to Lyapunov’s Central Limit Theorem $\frac{1}{\sqrt{\sum_{n=1}^N \sigma_n^2}} \sum_{n=1}^N \xi_n(z, t) \rightarrow \mathcal{N}(0, 1)$. Then, $\dot{z}_t = f(z, t) + g(z, t)b$ is an approximation of Stratonovich’s SDE (1), where $b \sim \mathcal{N}(0, 1)$. \square

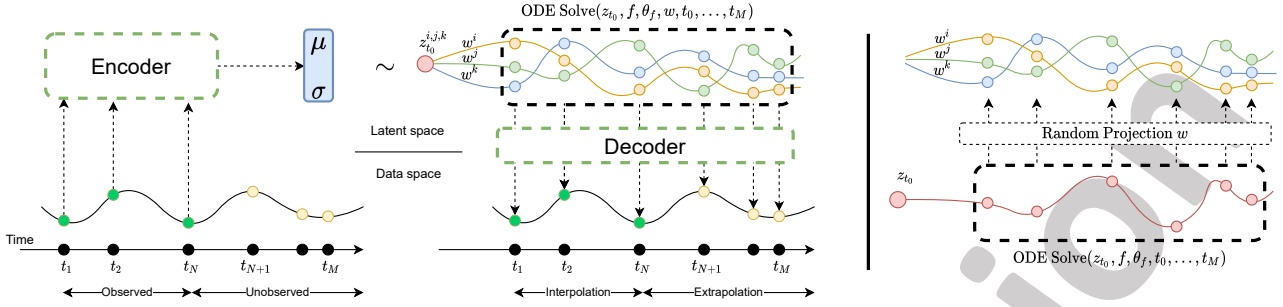


Figure 1: Structure of the model. First, encoder is applied to temporal data of an subject to generate initial point of the trajectory in latent space; Second, ME ODE solver is used to generate trajectory from the specified initial point; Last, decoder is used to map latent space ODE stages into observed values. On the right side we show how ME ODE can be viewed as random projection of trajectory of standard ODE, where trajectory is defined by random effect w .

This means that under standard second moment conditions on the stochastic part, the solution of (3) can be seen as an approximation of the solution to Stratonovich’s SDE in Equation (1). The benefits of this simplified form are twofold: **(a)** since z is a random variable, we can incorporate uncertainty within the ODE similar to a SDE, **(b)** for a given z , the trajectory in (3) can be modeled using an ODE and the associated computational benefits become available. Concurrently, [Hodgkinson et al., 2020] showed that it is possible to generalize this result to a more general class of SDEs using tools from the theory of rough paths. We must note (discussed briefly later), that for specific choice of f and g , the RHS of Equation (3) is well known in statistics and machine learning as a **Mixed Effects model** [Hyun et al., 2016]. Therefore, we will refer to our model as a mixed effects model and use the relevant terminology from this literature whenever possible. We hope that this will make the presentation more accessible and clarify that our scheme is *not a general purpose replacement to a deep neural network based SDE solver*.

Mixed effects model. Assuming individuals/groups are denoted by i , a nonlinear mixed effects (ME) model [Demidenko, 2013] can be written as:

$$\phi^i = \nu(X^i \beta + U^i \mathbf{b}^i) + \epsilon^i, \quad (6)$$

where $X^i \in \mathbf{R}^{n \times m}$ is a matrix of covariates where n and m are the number of observations and variables respectively. Here, ν is a non-linear (vector-valued) function, $\beta \in \mathbf{R}^m$ is a vector of *fixed* effects, $\mathbf{b}^i \sim \mathcal{N}(\mathbf{0}, \Sigma_b)$ is a vector of *random* effects, U^i is a design matrix (modeling choice) for random effects, $\phi^i \in \mathbf{R}^n$ is the response variable and $\epsilon^i \sim \mathcal{N}(\mathbf{0}, \Sigma_{\epsilon^i})$ represents a noise term.

Task. Our goal is to learn a latent representation of a time-varying physical process/dynamics z , and distribution $p(z|x_1, \dots, x_{t_k})$. Based on the Bayesian filtering discussion above, we focus on variational approximation techniques.

Learning latent representations with a VAE. Variational auto-encoders (VAE) [Kingma and Welling, 2013] enable learning a probability distribution on a latent space. Then, we can draw samples in the latent space – and the decoder can generate samples in the space of observations. In practice, the parameters of the latent distribution are learned by maximizing the *evidence lower bound* (ELBO) of the intractable likelihood:

$$\log p(\mathbf{x}) \geq -KL(q(\mathbf{z})||p(\mathbf{z})) + \mathbb{E}_{q(\mathbf{z})}[\log p(\mathbf{x} | \mathbf{z})] \quad (7)$$

where \mathbf{z} is a sample in the latent space from the approximate posterior distribution $q(\mathbf{z})$, with a prior $p(\mathbf{z})$, and \mathbf{x} is a reconstruction of a sample (e.g., an image) with the likelihood $p(\mathbf{x} | \mathbf{z})$. A common choice for q is $\mathcal{N}(\mu, \Sigma)$, where μ and Σ are trainable parameters [Kingma and Welling, 2013].

3 MIXED EFFECTS NEURAL ODE

Given a latent representation \mathbf{z} of a time-varying process with $\mathbf{z} = (\mathbf{z}_0, \dots, \mathbf{z}_n)$ and $\mathbf{z}_t \in \mathbf{R}^p$, we now model the latent representation \mathbf{z} as a mixed effect neural ODE. We will assume each of the p variables to be independent, so we will seek to learn p mixed effect models. Without any loss of generality, below we assume $z \in \mathbf{R}$ to denote the latent representation of the time-varying process at a time point t .

Modeling random effect \mathbf{b}^i in a network Γ . Mixed effects in the context of ODE is a well studied topic in longitudinal data analysis literature [Wang et al., 2014, Liang and Yu, 2013]. Formally, for subject i , given z^i , \mathbf{b}^i , and (population level) fixed effects β , we assume that there exist a smooth function h such that,

$$\dot{z}^i = h(z^i, \beta, \mathbf{b}^i). \quad (8)$$

Due to the universal approximation properties of neural networks [Zhou, 2020], such models are a sensible choice to express the nonlinear function h . Recall the non-linear mixed effects model from (6), and let us model

$$h(z^i, \beta, \mathbf{b}^i) = \nu(\eta(z^i)\beta + U^i \mathbf{b}^i).$$

Here, $\eta : \mathbf{R}^n \rightarrow \mathbf{R}^{n \times m}$ is a non-linear function, with n and m being the number of observations and variables respectively and $\beta \in \mathbf{R}^m$, $U^i \in \mathbf{R}^{n \times m}$. With a choice of $U^i = \eta(z^i)$, we can model $h(z^i, \beta, \mathbf{b}^i) = \Gamma(z^i)(\beta + \mathbf{b}^i)$, where Γ is a neural network with $\Gamma(z^i) \in \mathbf{R}^{n \times m}$. Now, we can derive the expressions for representing mixed effects in ODE, parameterized by a neural network as

$$h(z_t^i, \beta, \mathbf{b}^i) = \Gamma(z_t^i) \mathbf{w}^i, \quad (9)$$

where $\mathbf{w}^i \sim \mathcal{N}(\beta, \Sigma_b)$ is a mixed effect for subject i . This can be thought of as a projection from \mathbf{R}^m to \mathbf{R} along the direction given by $\mathbf{w}^i \in \mathbf{R}^m$.

Remark. Observe the difference between standard SDE in (1) – where the noise is added at each step t – and our formulation, where \mathbf{w}^i (or \mathbf{b}^i) is sampled once for subject i and completely defines the trajectory through $h(z^i, \beta, \mathbf{b}^i)$ for all steps of time t . This is *crucial* from the computational perspective: with this strategy, we can simply apply existing ODE solvers whereas backpropagating through a blackbox SDE requires specialized solutions, which are typically slower [Li et al., 2020, Liu et al., 2019].

Initializing ODE $h(z^i, \beta, \mathbf{b}^i)$ with an encoder E . Often in real-world analysis tasks involving panel data, the initial point of the process z_0 is not observed. While it can be learned as a parameter [Huang et al., 2008], it is desirable to also provide uncertainty pertaining to the learned z_0 . For this reason, we learn the distribution $q(z_0) = \mathcal{N}(\mu, \sigma)$, by training an encoder E to map observed data $\mathbf{x} = (\mathbf{x}_1, \dots, \mathbf{x}_n)$ (at all n time points) to parameters of $q(z_0)$, μ and σ :

$$(\mu, \sigma) = E(\mathbf{x}) \quad (10)$$

We use $q(z_0)$ to sample initial points of ODE z_0 .

Mapping z to x via decoder D . Given the latent representation z and the non-linear function D , which can recover the output \mathbf{x}_t^i , for a subject i at time point t , we can model the output of the dynamic process (e.g., in our application, a brain image) \mathbf{x}_t^i as a non-linear transformation of the latent measure of progression z_t^i :

$$\mathbf{x}_t^i = D(z_t^i) + \epsilon_t, \quad (11)$$

where ϵ_t is measurement error at each time point. This idea has been variously used in the literature [Pierson et al., 2019, Hyun et al., 2016, Whitaker et al., 2017].

The final model. Combining Equations (10), and (11) we obtain our Mixed Effects Neural ODE model in (12) and illustrated in Figure 1.

$$\begin{cases} z_0^i \sim \mathcal{N}(\mu, \sigma), \text{ where } \mu, \sigma = E(\mathbf{x}^i) \\ \mathbf{w}^i = \beta + b^i \sim \mathcal{N}(\beta, \Sigma_b) \\ \dot{z}_t^i = \Gamma(z_t^i) \mathbf{w}^i \\ \mathbf{x}_t^i = D(z_t^i) + \epsilon_t \end{cases} \quad (12)$$

Synopsis. Here, for subject i , we use the encoder E to map the observed data \mathbf{x}^i to parameters of the distribution of ODE initialization z_0^i . Then, we parameterize the derivative of ODE \dot{z}_t with a neural network $\Gamma(z_t^i)$ and mixed effects \mathbf{w}^i , and use D to map solution of ODE to the original space.

Structure of the latent space. When the latent space z can be embedded in a low dimensional space, it is natural to ask whether simulating the ODE \dot{z} in (12) can be accomplished efficiently. The following result shows a link between our model in (12) and random projection ideas [Vempala, 2005]. Moreover, in contrast to Neural ODE [Chen et al., 2018], random projections allow using high dimensional representations, $\Gamma(z_t^i)$, and mapping it back to \mathbf{R} using random projections. This provides expressive power but also approximately preserves the distance (using JL lemma).

Lemma 1 (Random projection). *With a certain choice of $h(z_t)$ in the ODE formulation and given a mixed effect \mathbf{w} with a choice of approximate posterior $q(\mathbf{w})$ as a Normal distribution, the solution to ME Neural ODE (Equation (12)) is a random projection of a solution to Neural ODE [Chen et al., 2018], Figure 1 (right part).*

Proof. Let us use the following notations,

- (a) $dz_t = f(z_t) dt$ defines trajectory of Neural ODE setup;
- (b) $d\tilde{z}_t = \Gamma(\tilde{z}_t) \cdot w dt$ defines trajectory of our ME setup;
- (c) $\tilde{z}_t = z_t \cdot w$: random projection of z_t . Then we have,

$$\begin{aligned} d\tilde{z}_t &= w \cdot dz_t = w \cdot f(z_t) dt = f\left(\frac{\tilde{z}_t}{w}\right) \cdot w dt \\ &= \Gamma(\tilde{z}_t) \cdot w dt \end{aligned}$$

If $\Gamma(\tilde{z}_t) = f\left(\frac{\tilde{z}_t}{w}\right)$, then \tilde{z}_t is random projection of z_t and $\tilde{z}_t = z_t \cdot w$. \square

While have a model, efficient training is still unresolved. Next, we show how for (12), we can derive ELBO-like bounds using Approximate Bayesian Computation (ABC) [Wilkinson, 2013, Fearnhead and Prangle, 2010].

Connection with approximation of Stratonovich’s SDE.

Observe that if we select f and g in (3) as $f(z, t) = \Gamma(z, t)\beta$ and $g(z, t) = \Gamma(z, t)\Sigma_b^{1/2}$, the approximation of Stratonovich’s SDE becomes ME-ODE defined in (9), i.e., we set $\dot{z}_t = \Gamma(z, t)\mathbf{w}$, where $\mathbf{w} = \beta + b\Sigma_b^{1/2} \sim \mathcal{N}(\beta, \Sigma_b)$.

Note that our ODE/SDE based derivation of the expression in (9) coincides with the mixed effects form proposed in [Xiong et al., 2019] for single panel (time) data.

Remark. Recall that in (8), we assumed that h is smooth. In theory, the smoothness assumption is justified due to Wong-Zakai approximation, see (3). In practice, this can be achieved by choosing a sufficiently fine discretization.

3.1 MODEL TRAINING: ME-NODE ELBO

The objective of the training scheme we describe now is to learn **(a)** distribution of z_0 , **(b)** fixed effect β , **(c)** variance of random effect Σ_b . To reduce clutter, in this section, we drop index i (which specifies a subject).

At a high level, our approach is to infer the random effects \mathbf{b} by learning it as a parameter. For our purposes, learning \mathbf{b} corresponds to ensuring that \mathbf{b} satisfies the following key requirement (accounting for a small reconstruction error): \mathbf{b} needs to be random by design for statistical reasons such as uncertainty quantification. Using our model in (12), it is easy to see that this requirement is satisfied because \mathbf{b} is sampled from $\mathcal{N}(\mathbf{0}, \Sigma_b)$. A common strategy to satisfy the requirement is to use a VAE [Chen et al., 2016]. It is known that in such probabilistic models computing the marginal likelihood $p(x)$ is usually intractable. Let $p(\mathbf{z}, \mathbf{w})$ be the prior joint distribution, $q(\mathbf{z}, \mathbf{w})$ as approximate joint posterior, and $p(x | \mathbf{z}, \mathbf{w})$ as likelihood of reconstruction. Using concepts from §2, we can derive a lower bound for the $p(x)$ of our ME-NODE model as:

$$\log p(x) = \log \int p(x | \mathbf{z}, \mathbf{w}) p(\mathbf{z}, \mathbf{w}) \frac{q(\mathbf{z}, \mathbf{w})}{q(\mathbf{z}, \mathbf{w})} d(\mathbf{z}, \mathbf{w}) \quad (13)$$

$$= \log \mathbb{E}_{q(\mathbf{z}, \mathbf{w})} \left(p(x | \mathbf{z}, \mathbf{w}) \cdot \frac{p(\mathbf{z}, \mathbf{w})}{q(\mathbf{z}, \mathbf{w})} \right) \geq \mathbb{E}_{q(\mathbf{z}, \mathbf{w})} \log p(x | \mathbf{z}, \mathbf{w}) - KL(q(\mathbf{z}, \mathbf{w}) || p(\mathbf{z}, \mathbf{w})), \quad (14)$$

where (13) follows from the marginalization property and then we use Jensen’s inequality. Next, we define $q(\mathbf{z}, \mathbf{w})$ which is needed to compute ELBO. Note that in the following description, $\xi | \psi$ refers to the random variable ξ conditioned on a value of ψ , regardless of what the value is, i.e., it can be $\xi | \psi = 0$ or $\xi | \psi = 1$.

Defining $q(\mathbf{z}, \mathbf{w})$. Assuming that z_0 and \mathbf{w} are independent random variables, we get

$$q(\mathbf{z}, \mathbf{w}) = q(z_0, \mathbf{z}_{-0}, \mathbf{w}) = q(\mathbf{z}_{-0} | z_0, \mathbf{w}) q(z_0) q(\mathbf{w}).$$

Recall that $\mathbf{z}_{-0} = (z_1, \dots, z_n)$ is a vector of ODE solutions at time step t , except $t = 0$. At every step t , z_t is a random variable, which is a function of z_0 and \mathbf{w} . However, with a fixed initial point z_0 and mixed effect \mathbf{w} , the progression follows a *defined* trajectory (i.e., there is no randomness). It means that $z_t | z_0, \mathbf{w}$ is deterministic and hence the distribution $q(z_t | z_0, \mathbf{w})$ is degenerate [Danielsson, 1994], which results in $q(\mathbf{z}_{-0} | z_0, \mathbf{w}) = \mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}}$. Thus,

$$q(\mathbf{z}, \mathbf{w}) = \mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}} q(z_0) q(\mathbf{w}) \quad (15)$$

Note. While the derivation from (13) to (14) is well defined for point mass distributions stated in (15), the use of ‘KL’, although consistent with [Bai et al., 2020] (pp 3, (4)–(6)) is not ideal (because $\log(0)$ and thus KL is undefined). We will avoid using KL notation in the loss in (16).

MC approximation of $E_{q(\mathbf{z}, \mathbf{w})} g(\mathbf{z}, \mathbf{w})$. The key in computing (14) is to estimate $E_{q(\mathbf{z}, \mathbf{w})} g(\mathbf{z}, \mathbf{w})$ for a given function $g(\mathbf{z}, \mathbf{w})$. Based on parameterization of $q(\mathbf{z}, \mathbf{w})$ in (15),

$$E_q g(\mathbf{z}, \mathbf{w}) = \int_{z_0, \mathbf{w}} g(\mathbf{z}, \mathbf{w}) \cdot \mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}} q(z_0) q(\mathbf{w}) dz_0 d\mathbf{w}.$$

While the integration may be intractable, it can be estimated by Monte Carlo (MC) techniques. Sampling (z_0^m, \mathbf{w}^m) from q , we compute $\frac{1}{M} \sum_{m=1}^M g^*(z_0^m, \mathbf{w}^m)$, where

$$g^*(z_0^m, \mathbf{w}^m) = g(z_0^m, \mathbf{w}^m) \cdot \mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0^m, \mathbf{w}^m \}}.$$

This type of sampling is called likelihood-free rejection sampling [Del Moral et al., 2012]: we reject all samples $(z_0^m$ and $\mathbf{w}^m)$, which do not generate observed $\mathbf{z}_{-0}^{\text{obs}}$.

The final loss. Given this MC approximation (with M samples), with the approximate posterior $q(\mathbf{z}, \mathbf{w})$ defined in (15) and with a similarly defined prior $p(\mathbf{z}, \mathbf{w}) = \mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}} p(z_0) p(\mathbf{w})$, the final loss is

$$\frac{1}{|S|} \sum_{s \in S} \left(\log p(x | z_0^s, \mathbf{w}^s) - \log \frac{q(z_0^s) q(\mathbf{w}^s)}{p(z_0^s) p(\mathbf{w}^s)} \right), \quad (16)$$

where S is a set: $\{ \forall s \in S : \mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0^s, \mathbf{w}^s \}} = 1 \}$.

Remark. While a MC approximation in (16) is an unbiased estimator of the Lower Bound in (14), its variance is $O\left(\frac{1}{|S|}\right)$. This leads to efficiency issues in that it may require a large M until we get z_0 and \mathbf{w} to generate \mathbf{z}_{-0} exactly along the observed trajectory to populate the set S . But we can address this problem using ABC methods [Wilkinson, 2013, Fearnhead and Prangle, 2010].

Efficient sampling: approximating $\mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}}$. ABC recommends finding samples of z_0 and \mathbf{w} to generate trajectories \mathbf{z}_{-0} which are *approximately* equal to the observed one, rather than exactly equal. The idea in [Marin et al., 2012] proposes using $\mathbb{1}_y \{ z \}$ as $\mathbb{1}_{A_{\epsilon, y}} \{ z \}$, where $A_{\epsilon, y} = \{ z | d\{z, y\} \leq \epsilon \}$ is an ϵ -neighborhood of y , and d is a distance function. For a direct application of these methods on $\mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}}$, we must have access to $\mathbf{z}_{-0}^{\text{obs}}$ in the latent space, which is unavailable unless the encoder E and the decoder D are identity functions. Nonetheless, we can approximate $\mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ \mathbf{z}_{-0} | z_0, \mathbf{w} \}}$, by comparing if the decoded $\mathbf{z}_{-0} | z_0, \mathbf{w}$ indeed corresponds to $\mathbf{z}_{-0}^{\text{obs}}$, i.e., we need to compute $\mathbb{1}_{\mathbf{z}_{-0}^{\text{obs}} \{ D(\mathbf{z}_{-0} | z_0, \mathbf{w}) \}}$. We simply use the mean squared error (MSE) as the distance for this comparison.

By decreasing ϵ , we can improve the quality of the samples for MC estimation, but at higher compute cost. However, because we learn the distribution of z_0 , during the first steps of training, our model provides poor reconstructions. For this reason, setting ϵ to a small value at the beginning of

the training is inefficient. Therefore, we make ε adaptive through the training, by choosing the sample, closest to our observed trajectory, i.e., sample with smallest distance d , and ε is a function of the initial point z_0 .

Choice of $q(\mathbf{w})$. To optimize the ELBO, it is necessary to define the approximate posterior $q(\mathbf{w})$ and prior $p(\mathbf{w})$. While we assumed in (9) and Lemma 1 that the true distribution of \mathbf{w} is Normal, $q(\mathbf{w})$ and $p(\mathbf{w})$ remain design choices for the user. For example, if we believe that the correlation structure of the data is sparse, then we have the following choices: Horseshoe [Carvalho et al., 2009], spike-and-slab with Laplacian spike [Deng et al., 2019] or Dirac spike [Bai et al., 2020]. However, for our experiments, we found that modeling $q(\mathbf{w})$ and $p(\mathbf{w})$ as Normal is sufficient.

Calibration. One feature of our model is that learned distribution of mixed effects \mathbf{w} can be used for personalized prediction during extrapolation [Wang et al., 2014, Ditlevsen and De Gaetano, 2005, Bouriaud et al., 2019]. First, we train our model to learn the parameters of distribution of z_0 , fixed effects β , and variance of random effect Σ_b . Then at test time, we make use of the observed temporal data for a previously unseen test subject \mathbf{x}^{obs} . Given the learned distribution of mixed effects \mathbf{w} , we want to find a sample $w \sim \mathbf{w}$, which minimizes error w.r.t. \mathbf{x}^{obs} . This selection provides the most appropriate mixed effect w corresponding to \mathbf{x}^{obs} . We call this process calibration: a solution to $\text{argmin}_{w \sim \mathbf{w}} \text{MSE}(\mathbf{x}^{\text{obs}}, \hat{\mathbf{x}}(w))$, where $\hat{\mathbf{x}}(w)$ is a prediction from our model. Note that this is slightly different from the average (used in probabilistic models like VAE).

Method summary. We provide a step-by-step summary,

Train and test phases

In the **training** phase, the observed data for subject i , x_t^i for $t \in [0, T]$ is assumed to be provided. Then, we

1. Use a suitable encoder E to map $\mathbf{x}^i = \{x_t^i\}$ to the latent representation of initial point z_0 of underlying ODE, as $z_0^i \sim \mathcal{N}(\mu, \sigma)$, where $\mu, \sigma = E(\mathbf{x}^i)$.
2. Given a suitable decoder D , we fit the ME-NODE model, by minimizing the loss in (16), thereby learning the appropriate distribution of mixed effects \mathbf{w}^i .

The output from this phase latent representation z_t^i described by ME-NODE model and the corresponding distribution of mixed effects \mathbf{w}^i .

In the **test** phase, the observed data for subject i , x_t^i for $t \in [0, T]$ is assumed to be provided. Then, we

1. Select a personalized mixed effect $w^i \sim \mathbf{w}^i$, according to the calibration scheme.
2. Then, we use the selected mixed effect sample w^i , to generate personalized prediction for the subject i for either interpolation or extrapolation.

The output from this phase is the prediction for subject i , \hat{x}_t^i for $t \in [0, \dots, T^*]$, where T^* can be extrapolated time, i.e., $T^* \geq T_i$, and/or a denser interpolation in $[0, T]$.

Estimated parameters	n_{z_0}, n_w				True values
	1, 1	1, 10	10, 1	10, 10	
$\hat{\mu}$	1.252	1.258	1.316	1.313	1.3
$\hat{\sigma}$	0.003	0.002	0.005	0.016	0.01
$\hat{\beta}$	0.311	0.315	0.311	0.319	0.3
$\hat{\sigma}_b$	0.036	0.051	0.054	0.060	0.01
MSE (all)	0.0017	0.0011	0.0006	0.0005	

Table 1: The first four rows show the estimated parameters for specific choices of n_{z_0} and n_w . Here we use the following convention: $(n_{z_0}, n_w) = (i, j)$ denotes we draw i and ij number of samples from z_0 and w respectively. The last row presents the MSE values for the estimated parameters.

4 EXPERIMENTS

We evaluate our model on five temporal datasets: **(1)** simulations, **(2)** MuJoCo hopper, **(3)** rotating MNIST, and **(4)** two different Neuroimaging datasets, representing disease progression in the brain.

Goals. We will evaluate: **(a)** ability to learn mixed effects, given different types of correlations in the data **(b)** the effect of mixed effects dimension m on extrapolation power and confidence of the model, and **(c)** the ability to preserve statistical group differences in the data in latent representations.

The baselines are given separately for each experiment. We provide description of hardware and neural networks architectures, including encoder/decoder in appendix.

4.1 SYNTHETIC DATASET

We start with a synthetic setup where all parameters are known. Using a ODE solver and conditioning on z_0 and w , we generate a solution of the mixed effect ODE

$$\begin{cases} z_0^i \sim \mathcal{N}(\mu = 1.3, \sigma = 0.01) \\ w^i \sim \mathcal{N}(\beta = 0.3, \sigma_b = 0.01) \\ \dot{z}_t^i = z_t^i w^i \end{cases} \quad (17)$$

We set the encoder E and decoder D in (12) to the identity transformation. Given 1000 (80 : 20 split for train/test) numerical solutions of the ODE in (17), we uniformly sampled 20 time points from $[0, 3]$, and use the first 10 time steps for interpolation and the last 10 for extrapolation.

Parameters. As the optimization of ELBO in (16) requires samples, we evaluate the performance of our model by varying the number of samples for z_0 and w (denoted by n_{z_0} and n_w respectively). The results in Table 1 suggest that MSE goes down with an increase of n_{z_0} or n_w .

In Figure 2 (*top panel*), we show samples (trajectories) drawn from the learned model (blue lines) with the real trajectories ('x' marker). Notice that the sampled trajectories from the learned model almost cover the "range" of real trajectories and the results appear meaningful.

Mixed effects. To evaluate generation of a personalized pre-

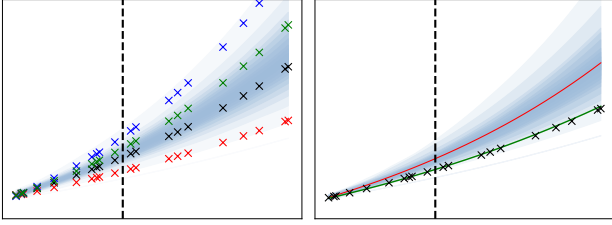


Figure 2: Real samples are denoted by “x” while blue lines are predictions. For each subfigure, the LHS and RHS of the dotted line contains interpolation and extrapolation results respectively. *Left*: the learned distribution of trajectories (from all test subjects) satisfies dynamic process described by the known system (1). *Right*: using the interpolation for calibration, we infer mixed effect w^i for i^{th} trajectory and generate a personalized prediction (green line). The prediction using BNN is shown in red.

diction for a subject i , by learning mixed effect w^i , recall that we split our data in two parts: interpolation (observed) and extrapolation (unknown). We use observed samples (interpolation part) to calibrate the mixed effect w^i and pass w^i to the selected trajectory during extrapolation. Figure 2 (*bottom panel*), shows that personalized prediction (green line) follows the observed data nicely across the entire time interval. In comparison, the standard BNN approach generates trajectory close to observed data for interpolation, and fails to extrapolate as well as our proposed model.

Runtime. For 1000 samples, runtime for an epoch of our method is 2.1 seconds, while Neural SDE takes about 27.13 seconds for the same memory utilization ($\sim 965\text{MB}$).

4.2 MUJOCO HOPPER

Here, we evaluate the performance of our model for simple Newtonian physics. Similar to NODE [Rubanova et al., 2019], we created a physical simulation using MuJoCo Hopper. While in [Rubanova et al., 2019], the generated samples were i.i.d, we explicitly introduce correlation between the samples. The process of MuJoCo Hopper is defined by the initial position and velocity. In order to generate correlated samples, we choose the initial velocity from the pre-specified set containing 1, 4 or 8 vectors. The entries of the velocity vectors are uniformly sampled from $[-2, 2]$. We evaluate our model on interpolation (10 steps) and extrapolation (10 steps) and compare results with NODE in Table 2. As our model implicitly learns correlation structure of the data by learning the distribution of mixed effect $q(w)$, we see an improvement in both interpolation and extrapolation. In addition, Figure 3 presents representative extrapolated samples using our proposed model.

4.3 ROTATING MNIST

We now evaluate a slightly more complicated rotating MNIST dataset. Here, we consider different types of correlations in the data and check: **(i)** relation between mixed effect dimension m (which we can think of as a dimension of random projection) and performance of the model, **(ii)** the



Figure 3: Visualization of 10 steps of extrapolation after observing 10 previous steps, with dimension of mixed effect $m = 50$. *Top* ground truth, *bottom* our prediction.

performance of personalized prediction for extrapolation in comparison with standard BNN.

Data description. Similar to the setup in ODE2VAE [Yildiz et al., 2019], we construct a dataset by rotating the images of different handwritten digits, in order to learn a digit specific mixed effect model. In ODE2VAE, digits were rotated by 22.5° . We used a slightly different scheme: for a sampled digit we randomly choose an angle from the set of 1, 4, or 8 angles from the range $[-\pi/4, \pi/4]$ and apply it at all time steps. For example, if we choose the set containing 4 angles, then a sampled digit is rotated using one of the 4 angles, selected randomly. In order to simulate a practical scenario, we spread out the initial points, by randomly rotating a digit by angles from $-\pi/2$ to $\pi/2$. We generate 10K samples of different rotating digits for 20 time steps and split it in two equal sets: interpolation and extrapolation.

Effect of mixed effect (random projection) dimension m .

Recall from Lemma 1, we showed that mixed effects w in (12) can be considered as a random projection. So, we can expect that with an increase in m , MSE should decrease as it leads to a richer latent representation of a trajectory. We see that this is indeed true as shown in Figure 4 (*left panel*). Here, we demonstrate the MSE of reconstruction for three values of m : 1, 20, and 50. We observe that for each time step, MSE decreases monotonically with an increase of m .

Calibration for personalized prediction. Here we choose $m = 50$ and compare our model for interpolation with ODE2VAE using 3 different rotation settings: 1, 4, and 8 available angles. The calibration results in Table 3 show that our model significantly outperforms the baseline ODE2VAE; however, making the correlation structure

		velocities		
		1	4	8
Interpolation	NODE	7.4	5.4	5.5
	This work	5.7	4.6	4.6
Extrapolation	NODE	166.1	82.1	80.3
	This work	164.1	81.2	80.0

Table 2: MSE (in scale of 10^{-3}) on MuJoCo Hopper data set, generated for three settings: 1, 4 and 8 initial velocities. We compare these two models using identical neural networks with the same number of levels and hyperparameters, however, in our model the dimension of mixed effect is $m = 50$.

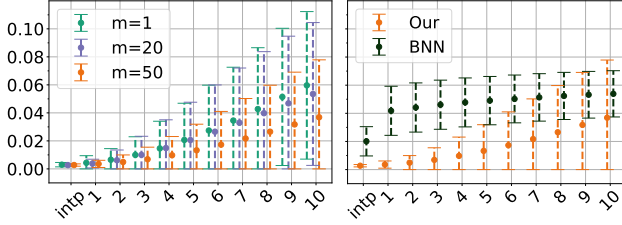


Figure 4: Distribution ($\mu \pm \sigma$) of MSE: *left*: varying with m : 1, 20, 50, *right*: first point on x-axis ('intp') indicates average MSE for all time steps: (1-10) of interpolation, and 1-10 indicate steps of extrapolation. Results for data with 8 possible angles shown.

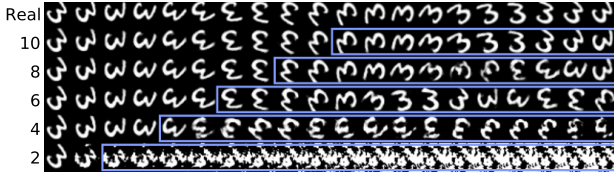


Figure 5: Visualization of extrapolated results (blue frame), given n time steps for calibration, where n is shown on y axis. The top row indicate real data.

of the data more complicated (increasing number of possible rotation angles), does lead to a larger MSE. This is expected: with an increase in complexity of correlation in data, the learning task (and thereby, prediction) becomes harder.

Recall that in order to generate a personalized prediction for subject i we have to sample mixed effect w^i resulting in trajectory closed to the observed. Thus, if the model fails to learn the distribution $q(w)$ accurately, sampling such w^i is less likely, and will result in a larger interpolation error.

For extrapolation, we compare with a standard BNN approach in Fig. 4. We observe that for each extrapolation step, we obtain, on average, much smaller MSE and smaller variance during the initial steps of extrapolation.

Extrapolation steps. Earlier, the number of steps for extrapolation were smaller than the number of observed steps, used for calibration. In Fig. 5, we show results of interpolation and extrapolation, varying the number of observed time steps used for calibration. Expectedly, decreasing the number of steps to be small for calibration yields a smaller number of steps where the extrapolation is meaningful.

		angles		
		1	4	8
Interpolation	ODE2VAE	0.0648	0.0644	0.0640
	Ours-50	0.0006	0.0014	0.0027

Table 3: MSE of two models, given different complexity of the data. Low interpolation error of our model indicates the properly learned mixed effect distribution $q(w)$.

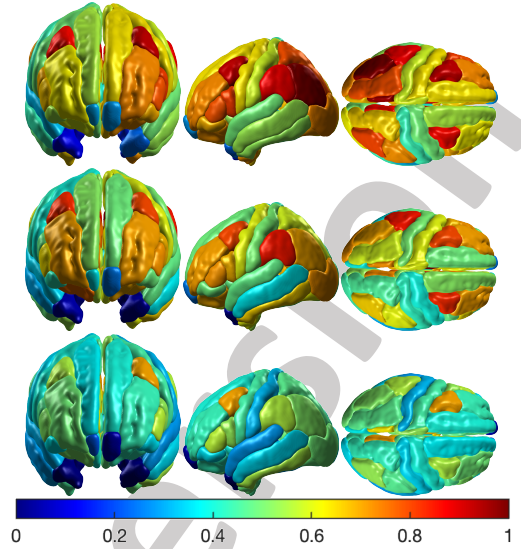


Figure 6: *Top - bottom row*: ground truth, our predicted, BNN predicted. We present the state of the brain at time step $t = 3$. Red and blue indicate high and low AV45 respectively. Compared to the baseline, our model is able to generate a sample for the subject with predicted values of AV45 closer to the ground truth.

4.4 LONGITUDINAL NEUROIMAGING DATA

In this section, we conduct experiments on two longitudinal brain imaging datasets obtained from Alzheimer’s Disease Neuroimaging Initiative (ADNI) (adni.loni.usc.edu), both of which describe AD progression through time, but are derived from two different imaging modalities.

Effect of number of dimensions m . We conducted experiments for different values of the mixed effect (random projection) dimension m , see appendix. We find that while for any m interpolation looks similar to real data, the further we move in extrapolation, the more noticeable the differences are. For example, for $m = 1$ some frames look blurry and in the last steps of extrapolation, the rotation is wrong. Increasing the dimension of random projection to $m = 20$ improves image quality, but does not fix rotation. Increasing dimension further to $m = 50$, not only improves quality of digits, but also leads to a better prediction of rotation.

(A) TADPOLE. TADPOLE dataset includes data for 276 participants with 3 time points. It represents Flortbetapir (AV45) Positron Emission Tomography (PET) scans, which measure the level of amyloid-beta pathology in the brain [Marinescu et al., 2018]. Scans were registered to a template (MNI152) to derive the 82 gray matter regions. Thus, each sample, at time t is a 82 dimensional vector, i.e., $x^t \in \mathbf{R}^{82}$.

Given a few time points (3 time points), we evaluate generation of a personalized prediction for a subject in an interpolation setting. We compare our personalized prediction with ground truth and prediction using a standard BNN approach. The predictions of both our model and BNN

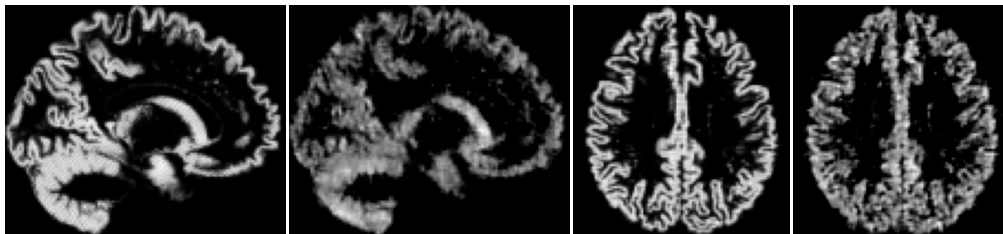


Figure 7: *Left-Right* Ground truth (1, 3) and prediction (2, 4) of two slices (side and top) of 3D brain at time point 3.

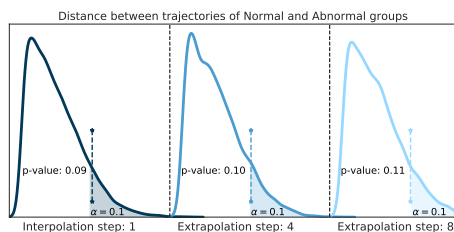


Figure 8: Distribution of distances resulted from permutation test with original distances (dashed lines), and corresponding p -values. The difference between trajectories is significant up to the 7th extrapolated step and degradation from 8.

are based on samples from the same distribution $q(z, w)$. However, Fig. 6 shows that the calibration of our model (Fig. 6, *second row*) provides better prediction than BNN (Fig. 6, *third row*). Even though the learned distribution of mixed effects $q(w)$ is capable of providing the correct trajectory (calibrated prediction), the direct application of the model without personalized calibration (BNN) leads to high subject-wise uncertainty.

(B) ADNI. Our second dataset from ADNI contains processed MRIs (3D brain scans) of size $105 \times 127 \times 105$ per subject at 3 time steps. The subjects are divided into two groups: diagnosed with Alzheimer’s disease (abnormal: 377 subjects) and healthy controls (normal: 152 subjects).

Given high resolution 3D images, we would like to evaluate whether our model is able to learn the distribution of mixed effects and perform calibration for personalized prediction. Similar to TADPOLE, we conduct an interpolation experiment and provide representative samples of brain images in Figure 7. We find by inspecting the axial/sagittal/coronal views that our model yields meaningful brain images. To evaluate the extrapolation capability, given limited number of time steps (only 3), we perform a statistical test. Recall that our method explicitly models the mixed effect term inside the trajectory to learn the data hierarchy. If our method works as intended, there should be a statistical difference between latent space of trajectories for normal and diseased/abnormal groups. Ideally, this difference should be preserved for several more extrapolation steps. To check this, we train our model on 3 time points. During testing, we use 3 observed time points for calibration and extrapolate for 5 more time steps: we get latent trajectories defined for 8 (3 interpolation and 5 extrapolation) time points. Finally, the resultant trajectories are used to evaluate differences between normal

and abnormal groups via a permutation test. The resultant distribution of distances and p -values for interpolation and extrapolation is in Fig. 8. As expected, for interpolation and some steps of extrapolation (up to step 7) differences between trajectories is significant (with p -value ≤ 0.1), and becomes less significant with more extrapolation steps.

5 DISCUSSIONS AND CONCLUSIONS

We proposed a novel ME-NODE model that enables us to incorporate both fixed and random effects for analyzing the dynamics of panel data. Our evaluations on several different tasks show that the ME-NODE loss function can be trained using existing ODE solvers in a stable and efficient manner. We see various benefits from incorporating mixed effects, **(1)** model explicitly learns the correlation structure of the data during the training, which improves prediction accuracy in setups where samples can be grouped by some criteria; **(2)** in contrast to generative models where only initial point is sampled from the distributions, by fixing initial point z_0 we can still provide uncertainty of the predictions, because from one initial point we can sample different trajectories; **(3)** since our model learns random effects for individual i , it allows personalized prediction, given a short history of data, which is useful in biomedical or scientific applications with a limited number of time points per individual along trajectory. One limitation of our approach is that there is not an explicit noise handling mechanism for test time calibration and prediction. This is problematic in large scale high dimensional settings. For example, say that the encoding distribution $\mathcal{N}(\mu, \Sigma)$ produces a small fraction of noisy trajectories. Even in small noise settings, filtering them for robust personalized prediction requires solving complex optimization problem [Bakshi and Kothari, 2021] and so handling noise is especially an open problem in real time, edge deployments. The code is available at https://github.com/vsingh-group/panel_me_ode.

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