# Manuscript progress report

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#### Method

#### Assumptions

- For each subject i in the population, a generalized outcome  $Y_i(t)$  is generated along a variable t (for example, time), where  $t \in (0, T)$ .
- The outcome, at any specific t, follows an exponential family distribution characterized by a (latent) continuous function  $\eta_i(t)$ :

$$g[E(Y_i(t))] = \eta_i(t) = \beta_0(t) + b_i(t)$$

$$p(Y_i(t)) = h(Y_i(t))exp\{\eta_i(t)T[Y_i(t)] - A(\eta_i(t))\}$$

• The continuous latent function consists of a population-level fixed process and an individual-level random process

$$\eta_i(t) = \beta_0(t) + b_i(t)$$

#### Observed data

In practice we would observe the discrete realization of  $\{Y_i(t), t\}$  along a dense grid. For simplicity, we assume the observation grid is regular (same across sample). When we have J observations points in (0, T], then for the jth observation point, we denote the corresponding value of t as  $t_j$ , and the corresponding outcome at this point  $Y_i(t_j)$ .

#### fGFPCA Algorithm

#### Bin data:

Choose a proper bin width w considering model complexity and identifiability. For now let's say the bins are equal-length and non-overlapping.

- Bin index s = 1...S
- Index of bin midpoints  $m_s$
- Value of t corresponding to bin midpoints  $t_{m_s}$
- Bin endpoints:  $(t_{m_s} \frac{w}{2}, t_{m_s} + \frac{w}{2}]$

#### Local GLMMs

At the every bin, we fit a local intercept-only model:

$$g[E(Y_i(t_j))] = \eta_i(t_{m_s}) = \beta_0(t_{m_s}) + b_i(t_{m_s})$$

where 
$$t_j \in (t_{m_s} - \frac{w}{2}, t_{m_s} + \frac{w}{2}].$$

Here we are basically saying that the value of latent function is constant within the same bin, which clearly is a misspecification of the true latent process.

From the model above, we will be able to estimate a  $\hat{\eta}_i(t_{m_s})$  on the binned grid for every individual in the training sample.

#### **FPCA**

Here, we fit a FPCA model on the  $\hat{\eta}_i(t_{m_s})$  obtained from step 2:

$$\hat{\eta}_i(t_{m_s}) = f_0(t_{m_s}) + \sum_{k=1}^K \xi_{ik} \phi_k(t_{m_s}) + \epsilon_i(t_{m_s})$$

where  $\xi_{ik}$  independently follows normal distribution  $N(0, \lambda_k)$ , and  $\epsilon_i(t_{m_s})$  at each point follows  $N(0, \sigma_2)$ . From this model, we will be able to obtain the following estimates which are shared across population:

- Population mean  $\hat{f}_0(t_{m_s})$
- Basis functions  $\hat{\mathbf{\Phi}} = \{\hat{\phi}_1(t_{m_s}), ..., \hat{\phi}_K(t_{m_s})\}$
- Estimates of variance of scores  $\hat{\lambda}_1...\hat{\lambda}_K$

#### **Projection and Debias**

The mean and basis functions are evaluated on the binned grid. To extend it to the original measurement grid data was collected on, we project the estimated eigenfunctions  $\hat{\Phi}$  back use spline basis. Now we have extend the  $\hat{\phi}_k(t_{m_s})$  to the original grid  $\hat{\phi}_k(t_j)$ 

Because of the misspecification of local GLMMs, the estimated eigenfunctions and eigenvalues are also biased by a constant multiplicative effect. Therefore, we use a GLMM to re-evaluate the mean function, eigenfunctions and eigenvalues.

#### Out-of-sample prediction

Now, let's assume we have a new subject u with  $J_u$  observations ( $J_u < J$ ). Then the log-likelihood of this new subject would be:

$$l_{u} = \sum_{t_{j} < t_{J_{u}}} log(h(Y_{u}(t_{j}))) + \hat{\eta}_{u}(t_{j})T(Y_{u}(t_{j})) - log(A[\hat{\eta}_{u}(t_{j})])$$

where 
$$\hat{\eta}_{u}(t_{j}) = \hat{f}_{0}(t_{j}) + \sum_{k=1}^{K} \xi_{uk} \hat{\phi}(t_{j})$$
.

With estimates for the population-level parameters from fGFPCA algorithms above, we can estimate  $\xi_{uk}$  by maximization of  $l_u$ . Direct maximization some times does not have closed form solution. Numeric maximization methods seem not very stable as well. So I have decided to used a Bayes approach (Laplace Approximation):

- Prior distribution:  $\xi_{uk} \sim N(0, \hat{\lambda}_k)$
- Posterior distribution: the likelihood of  $l_u = l(Y_u(t_i)|\xi_u)$

Laplace Approximation would get the posterior mode of  $\xi_{uk}$  through quadratic approximation.

### Larger-scale simulation

#### Simulation set up

Here we simulate binary data from cyclic latent process:

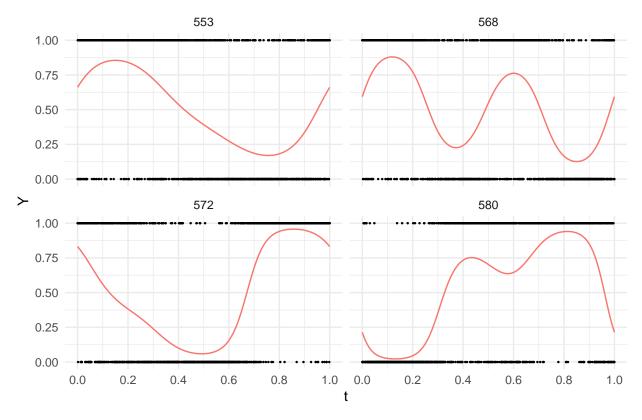
$$Y_{i}(t) \sim Bernoulli(\frac{exp(\eta_{i}(t))}{1 + exp(\eta_{i}(t))})$$

$$\eta_{i}(t) = f_{0}(t) + \xi_{i1}\sqrt{2}sin(2\pi t) + \xi_{i2}\sqrt{2}cos(2\pi t) + \xi_{i3}\sqrt{2}sin(4\pi t) + \xi_{i4}\sqrt{2}cos(4\pi t)$$

where:

- t is 1000 equal-spaced observations points on [0,1] (J = 1000).
- $f_0(t) = 0$
- $\xi_k \sim N(0, \lambda_k)$ , and  $\lambda_k = 1, 0.5, 0.25, 0.125$  for k = 1, 2, 3, 4 respectively.
- Sample size N = 500
- In the binning step, we bin every 10 observations
- 500 simulations were implemented

#### Simulated data



#### Reference method

- GLMMadaptive
- Here we can fit a model with random intercept and slope for time. It is doable on 500 datasets, but obviously too simple for the data generation scheme. We would expect it to perform terribly.

$$g(E(Y_i(t))) = \beta_0 + \beta_1 t + b_{i0} + b_{i1}t$$

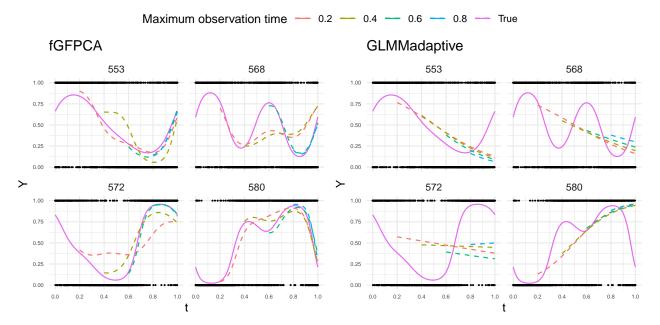
Table 1: Integrated squared error

	Maximum observation time								
		fGFF	PCA			GLMMa	daptive		
Window	0.2	0.4	0.6	0.8	0.2	0.4	0.6	0.8	
(0.2, 0.4]	146.407				387.708				
(0.4, 0.6]	183.967	74.977			291.579	269.799			
(0.6, 0.8]	218.265	49.275	15.776		315.778	282.736	278.242		
(0.8, 1.0]	108.918	77.981	17.747	12.005	563.011	477.485	597.746	600.34	

Table 2: Area under the ROC curve

	Maximum observation time								
	fGFPCA GLMMadaptive								
Window	0.2	0.4	0.6	0.8	0.2	0.4	0.6	0.8	
(0.2, 0.4]	0.748				0.591				
(0.4, 0.6]	0.664	0.734			0.524	0.596			
(0.6, 0.8]	0.715	0.790	0.803		0.669	0.694	0.687		
(0.8, 1.0]	0.740	0.755	0.781	0.784	0.514	0.556	0.526	0.564	

### Figure



#### **ISE**

#### **AUC**

I think we could say that while the total time spend on model fitting + prediction are similar between two methods, fGFPCA achieved much better flexibility and much better predictive performance of prediction

Table 3: Computation time (minutes)

Method	Fit	Prediction
fGFPCA GLMMadaptive	0.725 $2.287$	1.592 0.017

under every scenario.

#### Small-scale simulation: one iteration

Here we would like to fit fGFPCA and GLMMadaptive on a dataset with smaller sample size and/or smaller measurement density. For the GLMMadaptive model, we would set it up with spline basis functions so that its flexibility is comparable with fGFPCA model, such as:

$$g(E(Y_i(t))) = \sum_{k=1}^{5} \zeta_k B_k(t) + \sum_{l=1}^{5} \xi_{il} \phi_l(t)$$

I did the following things for data-reduction

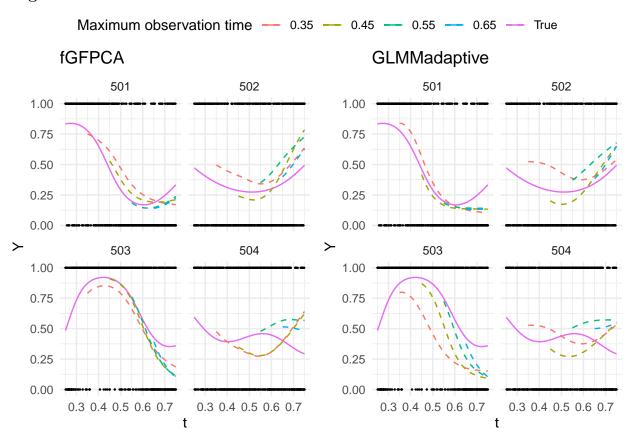
- Reduce sample size to 10 (for both training and testing)
- Reduce measurement track by half (0.25-0.75)
- Another way is the simple reduce the measurement density. For example, instead of every minute, we can use every 10 minutes. In this case, the binning step in fGFPCA would bin every 100 minutes in to one bin.
- — When we have smaller density of measurement grid, both methods could run into numeric issues in both methods. For example, singularity happened the the local GLMM step of fGFPCA procedure. This might be able to get resoved by changing the optimizer (e.g. Nealder\_Mead). Do we want to do that? Or perhaps we simply want to use the full track?
- We may also need to reduce number of simulations (GLMMadaptive took 20-ish minutes for one simulation).

Below are some of the outcomes:

Table 4: Integrated squared error

	Observation track									
		fGF1	PCA		GLMMadaptive					
Window	(0.25,0.35] $(0.25,0.45]$ $(0.25,0.55]$ $(0.25,0.65]$				(0.25, 0.35]	(0.25, 0.45]	(0.25, 0.55]	(0.25, 0.65]		
(0.35, 0.45]	35.724				48.779					
(0.45, 0.55]	77.379	22.690			74.537	34.232				
(0.55, 0.65]	32.581	20.370	20.043		48.683	35.666	18.303			
(0.65, 0.75]	158.191	135.968	130.360	66.648	174.158	141.285	143.370	103.92		

Figure



#### **ISE**

#### **AUC**

There is huge difference in computation time. GLMMadaptive took 23 minutes, while fGFPCA took less than 3 seconds.

### NHANES data application

We take 80% (7010) subjects for training, 20% (1753) for out-of-sample prediction.

Table 5: Integrated squared error

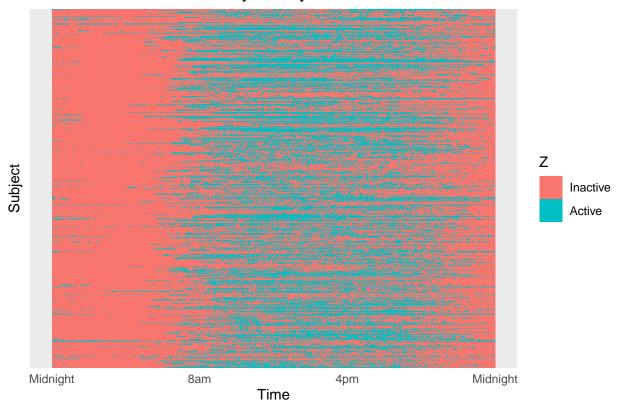
	Observation track									
		fGF1	PCA		GLMMadaptive					
Window	(0.25, 0.35]	(0.25, 0.45]	(0.25, 0.55]	(0.25, 0.65]	(0.25, 0.35]	(0.25, 0.45]	(0.25, 0.55]	(0.25, 0.65]		
(0.35, 0.45]	0.657				0.643					
(0.45, 0.55]	0.611	0.690			0.624	0.702				
(0.55, 0.65]	0.687	0.710	0.714		0.666	0.686	0.702			
(0.65, 0.75]	0.604	0.639	0.636	0.653	0.598	0.593	0.587	0.62		

Table 6: Area Under the ROC curve

	Max	Maximum observation time						
	fGF	PCA	GLMMa	adaptive				
Window	8am	4pm	8am	4pm				
8am-4pm 4am-midnight	$0.587 \\ 0.680$	0.766	0.628 0.448	0.613				

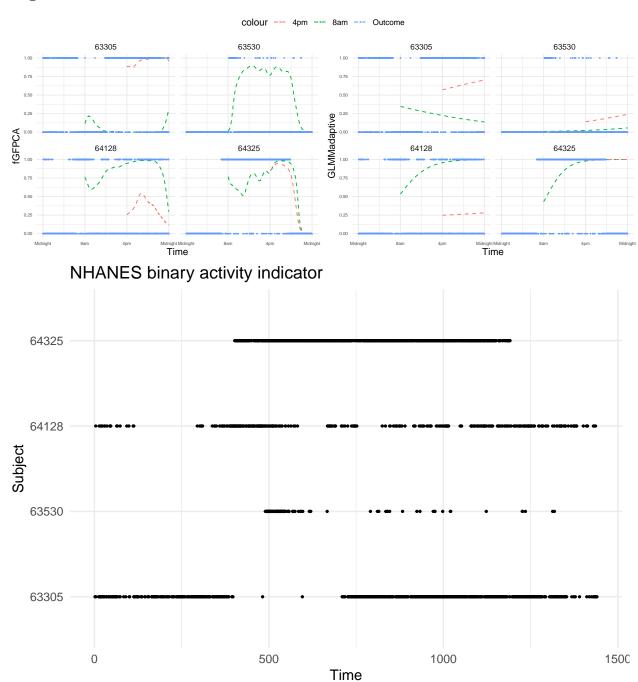
In the GLMMadaptive model, I could only fit a model with random intercept. It took more than 20 minutes for model fitting.

### Overview of NHANES binary activity indicator



AUC

## Figure



## Additional figures

### ${\it fGFPCA}$ algorithm demonstration

