

# An examination of the high-order dynamic interactions underlying multi-dimensional timeseries data

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## Abstract

Most complex systems reflect dynamic interactions between myriad evolving components (e.g., interacting molecules, interacting brain systems, interacting individuals within a social network or ecological system, coordinated components within a mechanical or digital device, etc.). Despite that these interactions are central to the full system’s behavior (e.g., removing a component from the full system can change the entire system’s behavior), dynamic interactions cannot typically be directly measured. Rather, the interactions must be inferred through their hypothesized role in guiding the dynamics of system components. Here we use a model-based approach to inferring dynamic interactions from timeseries data. In addition to examining first-order interactions (e.g., between pairs of components) we also examine higher-order interactions (e.g., that characterize mirrored structure in the patterns of interaction dynamics displayed by different subsets of components). We apply our approach to two datasets. First, we use a synthetic dataset, for which the underlying dynamic interactions are known, to show that our model recovers those ground-truth dynamic interactions. We also apply our model to a neuroimaging dataset and show that the high-order dynamic interactions exhibited by brain data vary meaningfully as a function of the cognitive “richness” of the stimulus people are experiencing.

## Introduction

The dynamics of the observable universe are meaningful in three respects. First, the behaviors of the *atomic units* that exhibit those dynamics are highly interrelated. The actions of one unit typically have implications for one or more other units. In other words, there is non-trivial *correlational structure* defining how different units interact with and relate to each other. Second, that correlational structure is *hierarchical* in the sense that it exists on many spatiotemporal scales. The way one group of units interacts may relate to how another group of units interact, and the interactions between those groups may exhibit some rich structure. Third, the structure at each level of this correlational hierarchy changes from moment to moment, reflecting the “behavior” of the full system.

These three properties (rich correlations, hierarchical organization, and dynamics) are major hallmarks of many complex systems. For example, within a single cell, the cellular components interact at many

spatiotemporal scales, and those interactions change according to what that single cell is doing. Within a single human brain, the individual neurons interact within each brain structure, and the structures interact to form complex networks. The interactions at each scale vary according to the functions our brains are carrying out. And within social groups, interactions at different scales (e.g., between individuals, family units, communities, etc.) vary over time according to changing goals and external constraints.

Although many systems exhibit rich dynamic correlations at many scales, a major challenge to studying such patterns is that typically neither the correlations nor the hierarchical organizations of those correlations may be directly observed. Rather, these fundamental properties must be inferred indirectly by examining the observable parts of the system— e.g., the behaviors of the individual atomic units of that system. In the *Methods* section, we propose a series of mathematical operations that may be used to recover dynamic correlations at a range of scales (i.e., orders of interaction). In the *Results* section, we demonstrate how our approach may be applied to multi-dimensional timeseries data: a synthetic dataset where the underlying dynamic correlations are known (we use this dataset to validate our approach), and a neuroimaging dataset comprising data collected as participants listened to a story (Simony et al., 2016). In different experimental conditions in the neuroimaging study, participants listened to altered versions of the story that varied in cognitive richness: the intact story (fully engaging), a scrambled version of the story where the paragraphs were presented in a randomized order (moderately engaging), a second scrambled condition where the words were presented in a random order (minimally engaging), and a “rest” condition where the participants did not listen to any version of the story (control condition). We use the neuroimaging dataset to examine how higher-order structure in brain data varies as a function of the cognitive richness of the stimulus.

## Methods

There are two basic steps to our approach. In the first step, we take a number-of-timepoints ( $T$ ) by number-of-features ( $F$ ) *matrix of observations* ( $\mathbf{X}$ ) and we return a  $T$  by  $\frac{F^2-F}{2}$  *matrix of dynamic correlations* ( $\mathbf{Y}$ ). Here  $\mathbf{Y}_0$  describes, at each moment, how all of the features (columns of  $\mathbf{X}$ ) are inferred to be interacting. (Since the interactions are assumed to be non-recurrent and symmetric, only the upper triangle of the full correlation matrix is computed.) In the second step, we project  $\mathbf{Y}_0$  onto an  $F$ -dimensional space, resulting in a new  $T$  by  $F$  matrix  $\mathbf{Y}_1$ . Note that  $\mathbf{Y}_1$  contains information about the correlation dynamics present in  $\mathbf{X}$ , but represented in a compressed number of dimensions. By repeatedly applying these two steps in sequence, we can examine and explore higher order dynamic correlations in  $\mathbf{X}$ .

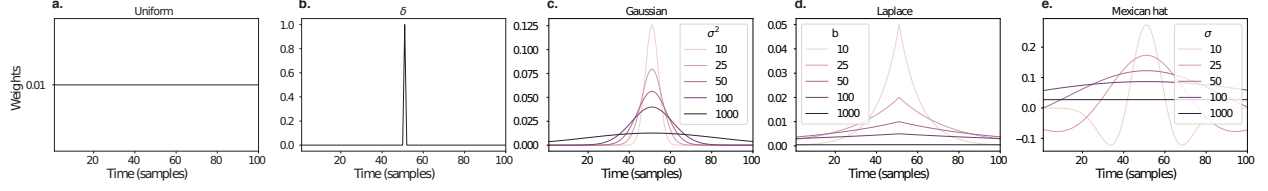


Figure 1: **Examples of time-varying weights.** Each panel displays per-timepoint weights at  $t = 50$ , evaluated for 100 timepoints (1, ..., 100). **a. Uniform weights.** The weights are timepoint-invariant; observations at all timepoints are weighted equally, and do not change as a function of  $t$ . This is a special case of weight function that reduces dynamic correlations to static correlations. **b. Dirac delta function.** Only the observation at timepoint  $t$  is given weight (of 1), and weights for observations at all other timepoints are set to 0. **c. Gaussian weights.** Each observation's weights fall off in time according to a Gaussian probability density function centered on  $\mu = t$ . Weights derived using several different example variance parameters ( $\sigma^2$ ) are displayed. **d. Laplace weights.** Each observation's weights fall off in time according to a Laplace probability density function centered on  $\mu = t$ . Weights derived using several different example scale parameters ( $b$ ) are displayed. **e. Mexican hat (Ricker wavelet) weights.** Each observation's weights fall off in time according to a Ricker wavelet centered on  $t$ . Weights derived using several different example width parameters ( $\sigma$ ) are displayed.

## 60 Dynamic correlations

Given a matrix of observations, we can compute the (static) correlations between any pair of observations,  $\mathbf{X}_i$  and  $\mathbf{X}_j$  using:

$$\text{corr}(\mathbf{X}_i, \mathbf{X}_j) = \frac{\sum_{t=1}^T (\mathbf{X}_i(t) - \bar{\mathbf{X}}_i)(\mathbf{X}_j(t) - \bar{\mathbf{X}}_j)}{\sqrt{\sum_{t=1}^T \sigma_{\mathbf{X}_i}^2 \sigma_{\mathbf{X}_j}^2}}, \text{ where} \quad (1)$$

$$\bar{\mathbf{X}}_k = \sum_{t=1}^T \mathbf{X}_k(t), \text{ and} \quad (2)$$

$$\sigma_{\mathbf{X}_k}^2 = \sum_{t=1}^T (\mathbf{X}_k - \bar{\mathbf{X}}_k)^2 \quad (3)$$

61 We can generalize this formula to compute time-varying correlations by incorporating a *weight function*  
 62 that takes a time  $t$  as input, and returns how much the observed data every timepoint (including  $t$ ) contribute  
 63 to the correlations at time  $t$  (Fig. 1).

Given a weight function  $w(t)$  for timepoint  $t$ , evaluated at timepoints in the interval  $[1, \dots, T]$ , we can

extend the static correlation formula in Equation 2 to reflect an *instantaneous correlation* at timepoint  $t$ :

$$\text{timecorr}(\mathbf{X}_i, \mathbf{X}_j, t) = \frac{\sum_{t=1}^T (\mathbf{X}_i(t) - \tilde{\mathbf{X}}_i(t)) (\mathbf{X}_j(t) - \tilde{\mathbf{X}}_j(t))}{\sqrt{\sum_{t=1}^T \tilde{\sigma}_{\mathbf{X}_i}^2(t) \tilde{\sigma}_{\mathbf{X}_j}^2(t)}}, \text{ where} \quad (4)$$

$$\tilde{\mathbf{X}}_k(t) = \sum_{i=1}^T w(t, i) \mathbf{X}_k(i), \quad (5)$$

$$\tilde{\sigma}_{\mathbf{X}_k}^2(t) = \sum_{i=1}^T (\mathbf{X}_k(i) - \tilde{\mathbf{X}}_k(t))^2, \quad (6)$$

and  $w(t, i)$  is shorthand for  $w(t)$  evaluated at timepoint  $i$ . Equation 5 may be used to estimate the instantaneous correlations between every pair of observations, at each timepoint (i.e.,  $\mathbf{Y}$ ).

## Higher-order correlations

Given a timeseries of dynamic correlations (e.g., obtained using Eqn. 5), higher-order correlations reflect the dynamic correlations between columns of  $\mathbf{Y}$ . Given unlimited computing resources, one could use repeated applications of Equation 5 to estimate these higher-order correlations (i.e., substituting in the previous output,  $\mathbf{Y}$ , for the input,  $\mathbf{X}$  in the equation). However, because each output  $\mathbf{Y}$  has  $O(F^2)$  columns relative to  $F$  columns in the input  $\mathbf{X}$ , the output of Equation 5 grows with the square of the number of repeated applications (total cost of computing  $n^{\text{th}}$  order correlations is  $O(F^{2(n-1)})$  for  $n \in \mathcal{J}, n > 0$ ). When  $F$  or  $n$  is large, this approach quickly becomes intractable.

To make progress in computing  $\mathbf{Y}_{n+1}$ , we can approximate  $\mathbf{Y}_n$  by computing an  $O(F)$ -dimensional embedding of  $\mathbf{Y}_n$ , termed  $\hat{\mathbf{Y}}_n$ , and then we can apply Equation 5 to  $\hat{\mathbf{Y}}_n$  rather than directly to  $\mathbf{Y}_n$ . This enables us to maintain  $O(n)$  scaling with respect to  $n$ , rather than exponential scaling via the direct approach.

There are many possible methods for computing  $\hat{\mathbf{Y}}_n$  from  $\mathbf{Y}_n$ , including traditional dimensionality reduction approaches and graph theory based approaches as described next. In the *Discussion* section we elaborate on other potential approaches.

## Dimensionality reduction-based approaches to computing $\hat{\mathbf{Y}}_n$

Commonly used dimensionality reduction algorithms include Principal Components Analysis (PCA; Pearson, 1901), Probabilistic PCA (PPCA; Tipping & Bishop, 1999), Exploratory Factor Analysis (EFA; Spearman, 1904), Independent Components Analysis (ICA; Jutten & Herault, 1991; Comon et al., 1991),  $t$ -Stochastic Neighbor Embedding ( $t$ -SNE; van der Maaten & Hinton, 2008), Uniform Manifold Approximation and Projection (UMAP; McInnes & Healy, 2018), non-negative matrix factorization (NMF; Lee & Seung,

1999), Topographic Factor Analysis (TFA) Manning et al. (2014), Hierarchical Topographic Factor analysis (HTFA) Manning et al. (2018), Topographic Latent Source Analysis (TLSA) Gershman et al. (2011), Dictionary learning (J. B. Mairal et al., 2009; J. Mairal et al., 2009), deep autoencoders (Hinton & Salakhutdinov, 2006), among others. While complete characterizations of each of these algorithms is beyond the scope of the present manuscript, the general intuition driving these approaches is to compute the  $\hat{\mathbf{Y}}$  with  $i$  columns that is closest to the original  $\mathbf{Y}$  with  $j$  columns, and where (typically)  $i \ll j$ . The different approaches place different constraints on what properties  $\hat{\mathbf{Y}}$  must satisfy and which aspects of the data are compared (and how) to characterize the match between  $\hat{\mathbf{Y}}$  and  $\mathbf{Y}$ .

Applying dimensionality reduction algorithms to  $\mathbf{Y}$  yields a  $\hat{\mathbf{Y}}$  whose columns reflect weighted combinations (or nonlinear transformations) of the original columns of  $\mathbf{Y}$ . This has two main consequences. First, with each repeated dimensionality reduction, the resulting  $\hat{\mathbf{Y}}_n$  has lower and lower fidelity (with respect to what the “true”  $\mathbf{Y}_n$  might have looked like without using dimensionality reduction to maintain scalability). In other words, computing  $\hat{\mathbf{Y}}_n$  is a lossy operation. Second, whereas the columns of  $\mathbf{Y}_n$  may be mapped directly onto pairs of columns of  $\mathbf{Y}_{n-1}$ , that mapping either becomes less cleanly defined in  $\hat{\mathbf{Y}}_n$  due to the reweightings and/or nonlinear transformations.

#### **Graph theory-based approaches to computing $\hat{\mathbf{Y}}_n$**

Graph theoretic measures take as input a matrix of interactions (e.g., using the above notation, an  $F \times F$  correlation matrix or binarized correlation matrix reconstituted from a single timepoint’s row of  $\mathbf{Y}$ ) and return as output a set of  $F$  measures describing how each node (feature) sits within that interactions matrix with respect to the rest of the population. Common measures include betweenness centrality (the proportion of shortest paths between each pair of nodes in the population that involves the given node in question; e.g., Newman, 2005; Opsahl et al., 2010; Barthélemy, 2004; Geisberger et al., 2008; Freeman, 1977); diversity and dissimilarity (characterizations of how differently connected a given node is from others in the population; e.g., Rao, 1982; Lin, 2009; Ricotta & Szeidl, 2006); Eigenvector centrality and pagerank centrality (measures of how influential a given node is within the broader network; e.g., Newman, 2008; Bonacich, 2007; Lohmann et al., 2010; Halu et al., 2013); transfer entropy and flow coefficients (a measure of how much information is flowing from a given node to other nodes in the network; e.g., Honey et al., 2007; Schreiber, 2000);  $k$ -coreness centrality (a measure of the connectivity of a node within its local sub-graph; e.g., Alvarez-Hamelin et al., 2005; Christakis & Fowler, 2010); within-module degree (a measure of how many connections a node has to its close neighbors in the network; e.g., Rubinov & Sporns, 2010); participation coefficient (a measure of the diversity of a node’s connections to different sub-graphs in the network; e.g., Rubinov & Sporns, 2010); and

sub-graph centrality (a measure of a node’s participation in all of the network’s sub-graphs; e.g., Estrada & Rodríguez-Velázquez, 2005).

As an alternative to the above dimensionality reduction approach to embedding  $\mathbf{Y}_n$  in a lower-dimensional space, but still allowing for scalable explorations of higher-order structure in the data, we also explore using the above graph theoretic measures as a means of obtaining  $\hat{\mathbf{Y}}_n$ . In particular: for a given graph theoretic measure,  $\eta : \mathcal{R}^{F \times F} \rightarrow \mathcal{R}^F$ , we can use  $\eta$  to transform each row of  $\mathbf{Y}_n$  in a way that characterizes the corresponding graph-theoretic properties of each column. Whereas the dimensionality reduction approach to computing  $\hat{\mathbf{Y}}_n$  is lossy, the graph-theory approach is lossless. However, whereas the dimensionality reduction approach maintains ties (direct or indirect) to the original activity patterns reflected in  $\mathbf{Y}_{n-1}$ , the graph-theory approach does not. Instead, the graph-theory characterizes the nature and timecourse of each feature’s *participation* in the network.

## Timepoint decoding

Cite HTFA paper, summarize relevant methods.

## Prediction

Use Eqn. 12 from SuperEEG paper, extend notation to treat  $\hat{\mathbf{Y}}_{1...n}(1...(t-1))$  as “observed” and use this to predict  $\mathbf{X}(t)$ . Key question: what is the predictive power added by each new level.

## Results

### Synthetic data

Neuroimaging dataset (Simony et al., 2016)

## Discussion

### Concluding remarks

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## Author contributions

Concept: J.R.M. Implementation: T.H.C., L.L.W.O., and J.R.M. Analyses: L.L.W.O and J.R.M.

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