1	Towards a psychology of individuals: the ergodicity information index and a bottom-up
2	approach for finding generalizations
3	Hudson Golino ¹ , John Nesselroade ¹ , & Alexander P. Christensen ²
4	¹ Department of Psychology, University of Virginia
5	2 Department of Psychology and Human Development, Peabody College, Vanderbilt
6	University

Author Note

Add complete departmental affiliations for each author here. Each new line herein
must be indented, like this line.

¹⁰ Enter author note here.

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The authors made the following contributions. Hudson Golino: Conceptualization,
 Methodology, Formal Analysis, Writing - Original Draft Preparation, Writing - Review &
 Editing; John Nesselroade: Writing - Original Draft Preparation, Writing - Review & Editing;
 Alexander P. Christensen: Writing - Original Draft Preparation, Writing - Review & Editing.
 Correspondence concerning this article should be addressed to Hudson Golino, 485
 McCormick Road, Gilmer Hall, Room 102, Charlottesville, VA 22903. E-mail:
 hfg9s@virginia.edu

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Abstract

In the last half of the 20th century, psychology and neuroscience have experienced a renewed 19 interest in intraindividual variation. To date, there are few quantitative methods to evaluate 20 whether a population structure (between-person) is likely to hold for individual people. We 21 present a network information theoretic approach to evaluate the extent to which a system 22 possesses the ergodic property. We introduce a new metric, the ergodicity information index 23 (EII), that can inform whether a set of multivariate time series (or a set of intensive 24 longitudinal measures) should be represented as multiple individual structures or as a single 25 population structure. The EII index quantifies the amount of information lost by 26 representing all individuals with a single population structure. If the individuals don't have a 27 similar structure, representing them with a single population network leads to a loss of 28 information. The EII index value will be higher than in cases where all individuals have a 29 similar structure. A Monte-Carlo simulation is implemented to test the EII index, and the 30 results show that the new index has a 94% accuracy in differentiating data in which all 31 individuals have a similar structure vs. data in which the individuals don't have a similar 32 structure. The paper also presents two new techniques designed to help applied researchers 33 to analyze data when the ergodicity property does not hold. The EII bootstrap test obtains 34 a sampling distribution of EII values as if all participants in the data have the population 35 network structure and compare this null distribution to the empirical EII value. Significant 36 differences indicate that the empirical data cannot be expected to be generated from an 37 ergodic process, and the population structure is not sufficient to describe all individuals. The 38 Domenico clustering method estimates the Von Neumann entropy of two networks and 39 computes their Jensen-Shannon Distance (JSD). Then, a complete-linkage agglomerative 40 hierarchical clustering method is applied to the JSD, and the clustering partitioning is 41 obtained via modularization maximization. The Domenico clustering method allows the 42 discovery of groups of individuals with a similar structure. Finally, two empirical examples 43 are shown, one using data from an intensive longitudinal experience sampling study 44

examining Big Five personality measured by the Big Five Inventory-2, and the other using 45 resting state neuroimaging data taken from a study examining creativity which used the 46 268-node Shen brain atlas. Starting with personality, the bootstrap EII test was significant, 47 suggesting that the BFI-2 data were nonergodic. Following up on the bootstrap EII test, the 48 information clustering was applied, and the single cluster test was performed. The single 49 cluster test suggested that the empirical networks had significantly greater JSD values than 50 the random networks meaning that the single cluster detected was not meaningful, and each 51 individual in the sample is unique. The brain networks had a significant bootstrap EII test 52 and significantly larger JSD values than random networks. These results, in line with the 53 personality data, suggest that (resting state) brain networks are not ergodic, and no 54 meaningful groups can be formed (i.e., each individual is unique). In sum, the personality 55 and (resting state) brain networks do not possess ergodicity and therefore lose information 56 when data are aggregated into a single population. 57

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Introduction

In the last half of the 20th century, psychology has experienced a renewed interest in 64 within-person (intraindividual) variation. One principal line of development sprang from the 65 introduction of P-technique factor analysis (Cattell, Cattell, & Rhymer, 1947). P-technique 66 involves the periodic collection (e.g. hourly, daily) of responses to a battery of measures 67 taken from a single person as opposed to across multiple people. The variation and 68 covariation of these measures are within a person (intraindividual) as opposed to the more 69 common between-person variation and covariation (interindividual) studied by differential 70 psychology. Although P-technique did not quickly become a standard tool for exploring 71 variability, its promise and the promise of studying intraindividual variation has attracted 72 increasing attention (Bereiter, 1963; Molenaar, 2004; Nesselroade & Ford, 1985). Today, 73 there is a considerable amount of research collecting data on intraindividual variability (Beck 74 & Jackson, 2022; e.g., ; Diehl, Hooker, & Sliwinski, 2014; Fisher, Reeves, Lawyer, Medaglia, 75 & Rubel, 2017; Gomes & Golino, 2015; Hultsch, Strauss, Hunter, & MacDonald, 2008; Ram, 76 Gerstorf, Lindenberger, & Smith, 2011), clearly having an impact on the field. 77

The re-emergence of within-person research, in an arena long dominated by between-person research, is not surprising. The accessibility of new technologies that are able to capture intraindividual variability (e.g., smartphones) has enabled researchers unparalleled opportunity to study a person rather than people. Intraindividual and interindividual variability has often been cast in terms of the idiographic versus nomothetic debate (see e.g., Lamiell, 1998). The recognition of these important thrusts has several consequences. There is an obvious need to formalize the definitions of both conceptions, and

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to do it in such a way that they can be readily distinguished at the conceptual level while
also highlighting their relation to each other (e.g., Molenaar, Huizenga, & Nesselroade, 2003;
Oertzen, Schmiedek, & Voelkle, 2020; Schmiedek, Lövdén, Oertzen, & Lindenberger, 2020;
Wright & Zimmermann, 2019).

Accompanying that effort, ways to distinguish between, and capitalize on the two kinds of variation in empirical data must also be created and made operational. For example, to what extent can we expect to characterize the way individuals differ from each other with the same structures that characterizes how a person changes over time? This question harks back to the old debate in developmental psychology about the relative merits of cross-sectional versus longitudinal research designs and whether cross-sectional data can be relied on to furnish accurate information regarding change over time at the individual level.

More recently, the ergodic property has surfaced in this context as a condition under 96 which one can expect within-person and between-person structures to match (Fisher, 97 Medaglia, & Jeronimus, 2018). If a system's process is ergodic, between-person structures 98 can be used to represent within-person structures (Molenaar et al., 2003) but, among 99 developmental psychologists at least, there seems to be little reason to think that ergodicity 100 will be a property of many (if any) developmental processes. To date, there are few 101 quantitative methods to evaluate whether a between-person structure is likely to hold for 102 individual people—that is, whether the system possesses the ergodic property. In the present 103 research, we present an information theoretic approach to evaluate the extent to which a 104 system possesses the ergodic property. From an information theoretic perspective, the 105 ergodic property can be framed in terms of the amount of information lost representing a set 106 of measures as a single between-person structure (nomothetic structure) instead of as 107 multiple within-person structures (within-person and idiographic structures). 108

Techniques to analyze complex systems with dynamic interactions between variables have a long history in statistical physics (e.g., Jaynes, 1957) and psychology (Boker, 2018;

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Cattell, 1965; Guttman, 1953), culminating in modern approaches such as network science 111 (Epskamp & Fried, 2018; Epskamp, Waldorp, Mõttus, & Borsboom, 2018a). In networks, 112 variables are represented as nodes (circles) and the edges (lines) between the nodes represent 113 associations between variables. There are many ways to characterize complex networks 114 (Newman, 2010), from the type of edge (e.g., directed or undirected, weighted or unweighted) 115 to the complexity of the network (e.g., algorithmic complexity; Morzy, Kajdanowicz, & 116 Kazienko, 2017; Zenil, Kiani, & Tegnér, 2018). Networks are a common method for 117 representing complex systems and the interactions between their components. 118

We introduce a new metric – termed *ergodicity information index* (EII) – that can 119 inform whether a set of variables should be represented as multiple individual structures 120 (multiple individual networks, multiplex networks) or as a single, between-individual 121 structure (unique single network, aggregate or population network). The EII characterizes 122 the relative algorithm complexity of the population structure with regard to multiple 123 individual networks taking into consideration the number of underlying dimensions (e.g., 124 communities, latent factors). Algorithm complexity of multiplex networks can be used to 125 determine the optimal number of layers needed to represent a multiplex network and to 126 detect structural and dynamical similarities among their layers (Santoro & Nicosia, 2020). 127 The representation of intraindividual structures as multiplex networks and the quantification 128 of their information relative to a single, population network structure are two central ideas in 129 the development of our EII. 130

The paper is organized as follows. The first section briefly introduces how intraindividual and interindividual structures can be estimated in a single framework using dynamic exploratory graph analysis (DynEGA; Golino, Christensen, Moulder, Kim, & Boker, 2022), an approach that combines generalized local linear approximation (Boker, Deboek, Edler, & Keel, 2010) and exploratory graph analysis (Golino & Epskamp, 2017; Golino, Shi, et al., 2020) to estimate dynamical factors in (intensive) longitudinal measures at different levels of analysis (individual, group, and/or population). In this section, we reframe the within- and between-person problem from a network perspective in which the intraindividual
structures are represented as a multiplex network (i.e., a collection of individual networks),
and the interindividual structure as a single (population) network.

In the second section, we develop the EII, introducing information theoretic concepts such as algorithm complexity that are necessary to define its meaning and interpretation. Next, a Monte Carlo simulation study is implemented to investigate the accuracy of the EII to differentiate between within- and between-person structures. Afterwards, we introduce an information theoretic approach to clustering, which can be used if the system is determined to be nonergodic.

Finally, one synthetic and two empirical examples from personality and neuroscience will be used to demonstrate how the EII can be used to determine whether a system possesses the ergodic property, and determine ergodic clusters (or groups) when the system is not. These new techniques are a step forward in the psychology of individuals, enabling the identification of generalizable constructs using a bottom-up approach (from individuals to group of individuals with common network characteristics).

¹⁵³ Representing intraindividual and interindividual structures as networks

To reframe the within- vs. between-person problem using a information theoretic 154 network approach, we first need to show how networks can be estimated in (intensive) 155 longitudinal data in a way that can generate both multiplex networks for the individuals 156 (i.e., multiple individual, within-person networks) and a single population or between-person 157 network. Recently, Golino et al. (2022) introduced a technique termed dynamic exploratory 158 graph analysis (**DynEGA**), combining techniques from dynamical systems (i.e., time-delay 159 embedding and generalized local linear approximation; Boker et al., 2010) and exploratory 160 graph analysis (Golino & Epskamp, 2017; Golino, Shi, et al., 2020) – a network psychometric 161 approach for dimensionality assessment and reduction. The DynEGA technique can be used 162

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to estimate dynamical communities (e.g., latent factors) in (intensive) longitudinal measures
at different levels of analysis (individual, group and/or population).

Network models have been proposed in psychological research for decades (e.g.,; Boker, 165 2018; Cattell, 1965; Guttman, 1953). More recently, a number of developments in network 166 modeling in psychology originated a new area termed *network psychometrics* (Epskamp, 167 2018) that relies mostly on the Gaussian graphical model (GGM; Lauritzen, 1996). 168 Assuming multivariate normality, a GGM can be obtained by modeling the inverse of the 169 variance-covariance matrix (and standardizing it to obtain partial correlations) in a way that 170 non-zero elements are freely estimated (Epskamp et al., 2018a), generating a sparse model of 171 the variance-covariance matrix (Epskamp, Rhemtulla, & Borsboom, 2017). As Epskamp et 172 al. (2018a) note, inverting and standardizing the variance-covariance matrix won't lead to 173 partial correlations that are exactly zero, meaning that the GGM is saturated. 174 Regularization techniques, such as a variant of the least absolute shrinkage and selection 175 operator (LASSO; Tibshirani, 1996) termed graphical LASSO (GLASSO; Friedman, Hastie, 176

¹⁷⁷ & Tibshirani, 2008), are generally used in network psychometrics (Epskamp & Fried, 2018;
¹⁷⁸ see; Epskamp et al., 2018a).

The GLASSO is a technique that is very fast to estimate both the model structure and the parameters of a sparse GGM (Epskamp, Waldorp, Mõttus, & Borsboom, 2018b). It has a tuning parameter (γ), that can be chosen in a way to minimize the extended Bayesian information criterion (EBIC; Chen & Chen, 2008), which is used to estimate optimal model fit and has been shown to accurately retrieve the true network structure in simulation studies (Epskamp & Fried, 2018; Foygel & Drton, 2010; Williams & Rast, 2020; Williams, Rhemtulla, Wysocki, & Rast, 2019).

The most used network estimation approach in psychological research (termed *EBICglasso*) is implemented in the **qgraph** package (version 1.4.1; Epskamp, Cramer, Waldorp, Schmittmann, & Borsboom, 2012). The **EBICglasso** function (Epskamp et al., ¹⁸⁹ 2012) samples 100 logarithmically-spaced values of λ , following Foygel and Drton (2010).

The ratio range of λ can be set by the user (defaults to 0.01). γ controls the severity of the model selection (defaults to 0.50). EBIC is computed for values of gamma larger than zero, and the value of λ that minimizes this information criteria is selected, generating a network with regularized edges.

Network psychometrics on cross-sectional data has progressed into dimensionality 194 assessment, where networks are estimated using the GGM or other network techniques (see 195 Golino, Shi, et al., 2020) and an algorithm for detecting communities in weighted networks 196 (Walktrap; Pons & Latapy, 2005) is used to identify latent factors. Golino and Epskamp 197 (2017) called this approach *exploratory graph analysis* (EGA). Simulation studies have found 198 EGA to perform as well as the most accurate factor analytic method, parallel analysis, and 199 produce the best large-sample properties of all the methods evaluated (Golino & Epskamp, 200 2017; Golino, Shi, et al., 2020). 201

In (intensive) longitudinal data, EGA can be used, but instead of using the 202 variance-covariance matrix of the raw data, it uses the variance-covariance of m-order 203 derivatives. The resulting network structure (GGM of the m-order derivatives) conveys 204 information on how variables are *changing* over time. As a result, the communities identified 205 using the Walktrap algorithm reflect not simple, static factors, but dynamical factors of 206 nodes that are fluctuating similarly as a function of time. Golino et al. (2022) called this 207 technique the dynamic exploratory graph analysis (DynEGA), merging network 208 psychometrics, dynamical systems modeling, and dimensionality assessment into a single 209 framework that can estimate structures at the individual or group/population levels. 210

DynEGA starts by transforming the time series of each variable $V = \{v_1, v_2, ..., v_N\}$ into a time delay embedding matrix $\mathbf{X}^{(n)}$, where *n* is the number of embedding dimensions. As pointed by Golino et al. (2022), a time delay embedding matrix is used to reconstruct the attractor of a dynamical system using a single sequence of observations (Takens, 1981; ²¹⁵ Whitney, 1936), preserving the phase-space dynamics of the system. In the time delay ²¹⁶ embedding matrix, each row is a phase-space vector (Rosenstein, Collins, & De Luca, 1993):

$$X = [X_1 \ X_2 \ \dots \ X_M]', \tag{1}$$

where X_i is the state of the system at discrete time *i* and is given by:

$$X_i = \begin{bmatrix} x_i \ x_{i+\tau} \ \dots \ x_{i+(n-1)\tau} \end{bmatrix},\tag{2}$$

where τ is the number of observations to offset successive embeddings (i.e., lag or reconstruction delay) and n is the embedding dimension. The time-delay embedding matrix is a $M \times n$ matrix, where $M = N - (n-1)\tau$ and N is the number of time points.

Once the time series of each variable collected in an (intensive) longitudinal study and transformed into a time-delay embedding matrix $\mathbf{X}^{(n)}$, derivatives can be estimated using generalized local linear approximation (Boker et al., 2010; **GLLA**; Deboeck, Montpetit, Bergeman, & Boker, 2009). Derivatives can represent different aspects of change such as the rate of change or velocity at which the variable is changing over time (first-order derivatives) and the speed of the rate of change or acceleration (second-order derivatives).

Deboeck et al. (2009) and Boker et al. (2010) show how derivatives can be estimated in the GLLA framework:

$$\mathbf{Y} = \mathbf{X}\mathbf{L}(\mathbf{L}'\mathbf{L})^{-1},\tag{3}$$

where **Y** is a matrix of derivative estimates, **X** is a time delay embedding matrix (with *n* embedding dimensions; to simplify the notation, $\mathbf{X} = \mathbf{X}^{(n)}$), and **L** is a matrix with the weights expressing the relationship between the embedding matrix and the derivative estimates. The weight matrix **L** is a $n \times \alpha$ matrix, where *n* is the number of embedding dimensions and α is the (maximum) order of the derivative. Each column of the weight matrix is estimated as follows, considering the order of the derivatives going from zero to k, $\alpha = [0, 1, ..., k]$:

$$\mathbf{L}_{\alpha} = \frac{\left[\Delta_t (v - \bar{v})\right]^{\alpha}}{\alpha!} \tag{4}$$

where Δ_t is the time between successive observations in the time series, v is a vector from one to the number of embedded dimensions (i.e., v = [1, 2, ...n]), \bar{v} is the mean of v, α is the order of the derivative of interest, and α ! is the factorial of α .

After estimating the derivatives for all time series (i.e., all variables). EGA is used to 239 estimate the multiplex networks (intraindividual or within-person structures). In this 240 process, each matrix of derivatives will generate a different network and dimensionality 241 structure (represented by clusters of nodes in the network) for each individual. The 242 population (between-person or nomothetic) structure can be estimated by stacking the 243 derivative matrix of each individual (i.e., row-binding the matrices) and applying EGA to the 244 stacked matrix. In both cases (i.e., individual or population structures), the resulting clusters 245 in the networks corresponds to variables that are changing together (Golino et al., 2022). 246

An example illustrates how this technique works: Suppose we ask two people to answer 247 eight items of depression once a day for 100 days. After applying DynEGA to the data, we 248 would obtain three important types of information (see Figure 1): the derivatives for the 249 eight variables, the network structure for each person (intraindividual), and the network 250 structure of the two people combined (interindividual). The left side of Figure 1 shows the 251 time series of first-order derivatives while the network structure of each individual are at the 252 center (top and bottom represents subject one and two, respectively). The interindividual 253 network structure is depicted in the right side of Figure 1. As can be seen in the network 254 structure of person one (top network at the center of Figure 1), the pair of nodes pink and 255 green are not connected (i.e. there is no edge linking these pairs of nodes). However, the 256

same pair of nodes are connected in the network structure of person two (bottom network at 257 the center of Figure 1), and in the between-individual network structure (or population 258 network) depicted at the right side of Figure 1. A question that must be answered now is: 259 how much information is being lost by representing the structure of the two individuals as a 260 single, interindividual network, relative to representing them as a multiplex network? Can 261 the derivatives of person one and two be stacked to estimate a single network or by doing so, 262 is important information about each individual lost? The next section aims to answer these 263 questions and to introduce the ergodicity information index. 264





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The Ergodicity Information Index

The individual networks and population network in Figure 1 can be compared in terms of their algorithm complexity. Algorithm complexity can be used to analyze complex objects in an unbiased manner using mathematical principles (Zenil et al., 2018), and is based on the work of Kolmogorov (1968), Martin-Löf (1966), Solomonoff (1964) and others. As Zenil et al. (2018) and Morzy et al. (2017) show, the algorithm (or Kolmogorov) complexity of a string sis formally defined as:

$$K_s = \min(|P|, T(P) = s)$$

where P is a program producing the string s when running on a universal Turing 272 machine T, and |P| is the number of bits required to represent P (i.e., the length of P). A 273 Turing machine is a formal model of a general-purpose computer that can be programmed to 274 reproduce any computable object, such as a string (Zenil et al., 2018). The Kolmogorov 275 complexity of a string is defined, in other words, as the length of the shortest possible 276 program that can produce that string as its output (Santoro & Nicosia, 2020). Kolmogorov 277 complexity has a drawback of being incomputable (Zenil et al., 2018), since there is no way 278 to estimate the number of possible programs that could produce the string s (Kolmogorov, 279 1968). But it can be approximated by using compression algorithms (Morzy et al., 2017; 280 Santoro & Nicosia, 2020), in which the compressed string s is an estimate of K_s . 281

To obtain an estimation of Kolmogorov complexity in networks, the most common 282 approach is to compute the size of the compressed weighted edge list (Santoro & Nicosia, 283 2020). For single, unique networks, this is straightforward, but for multiplex networks, the 284 Kolmogorov complexity requires a strategy to encode all individual graphs into a single 285 network. Santoro and Nicosia (2020) proposed the use of a prime-weight encoding matrix Ω 286 that assigns a distinct prime number $(p^{[\alpha]})$ to each individual network (i.e., each of the A 287 layers of the multiplex networks) and sets each element Ω_{ij} equal to the product of the 288 primes associated to the layers where an edge between node i and j exists: 289

$$\Omega_{ij} = \begin{cases} \prod_{\alpha:\alpha_{ij}^{[\alpha]}=1} p^{[\alpha]} \\ 0 \text{ if } \alpha_{ij}^{[\alpha]} = 0 \ \forall \ \alpha = 1, \dots, A \end{cases}$$
(5)





Santoro and Nicosia (2020) proposed a new metric for quantifying the algorithm complexity of multiplex networks that can be computed as the ratio of the (approximate) Kolmogorov complexity of the prime-weight matrix Ω of a multiplex network with A layers and the Kolmogorov complexity of an aggregated network combining all layers.

A similar strategy can be followed to quantify the algorithm complexity of the networks estimated using DynEGA. The multiplex networks are all individual networks estimated using the derivatives computed via generalized local linear approximation in the DynEGA technique. Instead of comparing the algorithm complexity of Ω with an weight aggregation of the multiplex networks, it is more informative to compare it with the 299

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³⁰¹ also important to consider the number of latent factors underlying the intensive-longitudinal
³⁰² data. Therefore, our *ergodicity information index* can be computed as:

$$\xi = \sqrt{F_P} \begin{bmatrix} \frac{\binom{(K_{\Omega})}{K_{P*}}}{\log(L_{\chi})} \end{bmatrix}$$

where $\sqrt{F_P}$ is the square-root of the number of factors estimated in the population structure using dynamic exploratory graph analysis, K_{Ω} is the algorithm complexity of the prime-weight encoding matrix of the individual networks (that composes the multiplex network χ), K_{P*} is the algorithm complexity of the prime-weight transformation of the population network (i.e., each element in the population network, P_{ij} , is transformed such that $P_{ij}*=2^{P_{ij}}$), and $L\chi$ is the number of distinct edges across the networks that make up the multiplex network (i.e., non-zero edges).

The EII (ξ) computes the amount of information lost representing a set of measures as a single interindividual structure (nomothetic structure) instead of representing the measures as multiple individual structures (within-person or intraindividual structures). Larger values of the ergodicity information index indicate that the intraindividual networks encode a relatively larger amount of information with respect to the population network.

The ratio $\frac{(K_{P*})}{K_{\Omega}}$ is normalized by the number of distinct edges in the multiplex network (set of individual networks) because a set of networks with larger number of edges is expected to have a higher complexity. Similar to Santoro and Nicosia (2020), we implemented a canonical prime association to compute ξ . In canonical prime association, prime numbers are associated to layers in increasing order of their total number of edges.

In terms of estimation, the Kolmogorov complexity is estimated for each network (i.e., prime-weight transformation of the population network P* and the prime-weight encoding of

the multiplex networks Ω) as the length of the compression of the string formed by their 322 edge list, using the *Gzip* compression algorithm available in the memCompress function of 323 base R (R Core Team, 2017). In theory, Kolmogorov complexity can also be computed using 324 other representations of the network such as a string composed by the rows of the Laplacian 325 matrix, degree list, degree distribution, or weights list (Morzy et al., 2017). In this paper, we 326 focus mainly on the algorithm complexity estimated using the weights list, and compare it to 327 the estimation using the edge list. Since Kolmogorov complexity is affected by the order of 328 elements in the edge list, the final estimation of algorithm complexity is based on the mean 329 of 1,000 computations of K over an edge list randomly ordered. One advantage of using 330 Kolmogorov complexity to estimate the complexity of networks is that it is less dependent on 331 the network representation than other metrics of complexity such as entropy-based metrics 332 (Morzy et al., 2017). 333

The use of the EII implies a different type of ergodicity that we call *super-weak* 334 ergodicity. In a strict definition of ergodicity, there are two central requirements: stationarity 335 for all participants and homogeneity for all time points (Voelkle, Brose, Schmiedek, & 336 Lindenberger, 2014). A softer type of ergodicity termed *weak ergodicity*, requires only that 337 the marginal distributions for all participants and for all time points be identical (Oertzen et 338 al., 2020). The super-weak ergodicity, on the other side, doesn't require stationarity for all 339 participants (i.e., the same covariance matrix for all subjects), homogeneity for all time 340 points (i.e., the same covariance matrix across time), or an equal marginal distributions for 341 all participants and time points. It requires a much weaker condition: the algorithm 342 complexity of the population (or between-person) network be similar (but not equal) to the 343 algorithm complexity of the prime-weight encoded network of all individuals. 344

Suppose we have four people that are assessed using an eight-item questionnaire for 100 days. Persons one and two are more similar than persons three and four, although none of them are exactly equal to one another. DynEGA is used to compute the first-order derivatives for each variable and to generate a network for each individual, and two ³⁴⁹ between-person or population networks: one for persons one and two, and one for persons ³⁵⁰ three and four. If we calculate the correlation of the derivatives for each variable, none of ³⁵¹ them are exactly equal to the others. Figure 3 shows the heatmap of the correlation matrices ³⁵² (calculated using the first-order derivatives of each variable) and the resulting network ³⁵³ structure for each person and each population with node colors representing the estimated ³⁵⁴ latent factors.

Two factors were identified in individuals one, two and three, and four factors were 355 identified in individual four. The population network for persons one and two indeed shows 356 two factors, while the population network for persons three and four shows four factors, both 357 estimated using DynEGA. Calculating the ergodicity information index (ξ), we obtain $\xi =$ 358 1.14 for individuals one and two and $\xi = 1.29$ for individuals three and four. Therefore, more 359 information is lost by representing the eight measures as a single structure for individuals 360 three and four (bottom of Figure 3, population 2) than for individuals one and two (top of 361 Figure 3, population 1). Another way to interpret the results above is that the 362 intraindividual networks for individuals three and four encode a relatively larger amount of 363 information with respect to the population network, compared to individuals one and two 364 and their population structure. 365

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Simulation Design

To verify the suitability of the ergodicity information index to identify if intensive 367 longitudinal measures should be represented as a set of within-person structures (multiple 368 individual networks) or as a single, between-person or population structure (only one 369 network), a Monte Carlo simulation is implemented. Four data conditions were 370 systematically manipulated: sample size (10, 50 and 100), number of time points (50 and 371 100), number of variables per factor (4 and 6) and number of factors (2, 3). Two separate set 372 of conditions were used in the simulation. In the first, all individuals had the same number 373 of factors (Eq condition). Therefore, representing these individuals using a single population 374



Figure 3

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structure is reasonable. In the second set of conditions, half of the subjects had the same
number of factors (and variables per factor) as condition one, but the other half had a
different configuration, with more or less factors (*NotEq* condition).

The sample sizes were selected to reflect low, moderate, and high samples, consistent 378 with many empirical papers using an intensive longitudinal measurement design that 379 typically don't use data from more than 100 individuals (Liu, Zhou, Palumbo, & Wang, 2016; 380 Schmiedek, Lövdén, & Lindenberger, 2020). Regarding the number of time points and factor 381 loadings (set to unity), they were selected based on the conditions tested by Zhang, 382 Hamaker, and Nesselroade (2008). In terms of the number of variables per factor, three are 383 the minimum required for factor identification (Anderson, 1958). In the present simulation, 384 the number of items were selected to reflect adequate (4) and slightly overindentified (6)385 factors (Velicer, 1976; Widaman, 1993). The measurement error covariance matrix had 0.1 in 386 the diagonal, so we could see how the EII works under minimum error conditions. This 387 strategy allows for the impact of the other variables systematically manipulated in the 388 simulation to be better understood. 389

The autoregression coefficients of the DFAS were set to 0.1, 0.4, and 0.8, representing a low, moderate, and high autoregressions. Additionally, two conditions were held constant: the matrix with autoregressive (0.8) and cross-regressive coefficients (0), and the covariance matrix for the random shock (off-diagonal = 0.18; diagonal = 0.36). The values of the cross-regressive coefficients and the random shock matrix were selected following Zhang et al. (2008).

396 Data Generation

In the Monte Carlo simulation, 500 data matrices were generated for each combination of variables (number of factors, number of variables per factor, number of time points, sample size, autoregression coefficients) according to the DAFS model. First, the matrix of

random shock vectors \mathbf{v}_t was generated following a multivariate normal distribution with 400 mean zeros and $q \times q$ covariance matrix **D** (off-diagonal values = 0.18; diagonal values = 401 (0.36), where q is the number of factors and t is the number of time points plus 1,000 (used as 402 the burn-in estimates for the Markov chain). Second, the factor scores are calculated and the 403 first 1,000 estimates are removed (burn-in phase). Third, the measurement error matrix is 404 estimated following a multivariate normal distribution with mean zeros and $p \times p$ covariance 405 matrix Q, where p is the total number of variables (number of variables per factor times F). 406 Finally, the observed variables \mathbf{Obs}_t at time $t \ (t = 1, 2, ..., N)$ are calculated using the 407 following equation 6. 408

$$\mathbf{Obs}_t = \mathbf{\Lambda} \mathbf{F}_t + \mathbf{e}_t,\tag{6}$$

where Λ is the factor loading matrix $(p \times q)$, \mathbf{F}_t is a $q \times 1$ vector of factors at time t, and \mathbf{e}_t is a $p \times 1$ vector with measurement errors following a multivariate normal distribution with mean zeros and covariance matrix Q (Nesselroade, McArdle, Aggen, & Meyers, 2002; Zhang et al., 2008).

413 The factor scores, \mathbf{F}_t , are calculated as follows:

$$\mathbf{F}_{t} = \sum_{l=1}^{L} \mathbf{B}_{l} \mathbf{F}_{t-l} + \mathbf{v}_{t}$$
(7)

where \mathbf{B}_l is a $q \times q$ matrix of autoregressive and cross-regressive coefficients, \mathbf{F}_{t-l} is a vector of factor score l occasions prior to occasion t and \mathbf{v}_t is a random shock vector (or innovation vector) following a multivariate normal distribution with mean zeros and $q \times q$ covariance matrix \mathbf{D} (Nesselroade et al., 2002; Zhang et al., 2008). In the DAFS model, $\mathbf{\Lambda}$, \mathbf{B}_l , \mathbf{Q} and \mathbf{D} are invariant over time.

Data following the DAFS model can be simulated using the simDFM function of the *EGAnet* package (version 1.2.0; Golino & Christensen, 2019).

Results

Table 1Accuracy per conditions tested

Ν	TimeP	NFAC	NVAR	NFAC2	NVAR2	Autoreg	Accuracy	Lower 95% C.I	Upper 95% C.I
10	50	2	4	4	2	0.1	75.00	72.31	77.69
10	50	2	4	4	2	0.4	91.30	89.55	93.05
10	50	2	4	4	2	0.8	98.70	98.00	99.40
10	50	2	6	3	4	0.1	100.00	100.00	100.00
10	50	2	6	3	4	0.4	100.00	100.00	100.00
10	50	2	6	3	4	0.8	100.00	100.00	100.00
10	50	3	4	2	6	0.1	74.40	71.69	77.11
10	50	3	4	2	6	0.4	68.20	65.31	71.09
10	50	3	4	2	6	0.8	68.90	66.03	71.77
10	50	3	6	2	9	0.1	78.50	75.95	81.05
10	50	3	6	2	9	0.4	90.30	88.46	92.14
10	50	3	6	2	9	0.8	98.80	98.12	99.48
10	100	2	4	4	2	0.1	70.00	67.15	72.85
10	100	2	4	4	2	0.4	93.50	91.97	95.03
10	100	2	4	4	2	0.8	99.70	99.36	100.04
10	100	2	6	3	4	0.1	100.00	100.00	100.00
10	100	2	6	3	4	0.4	100.00	100.00	100.00
10	100	2	6	3	4	0.8	100.00	100.00	100.00
10	100	3	4	2	6	0.1	83.60	81.30	85.90
10	100	3	4	2	6	0.4	83.20	80.88	85.52
10	100	3	4	2	6	0.8	83.78	81.49	86.07
10	100	3	6	2	9	0.1	82.40	80.04	84.76
10	100	3	6	2	9	0.4	96.90	95.82	97.98
10	100	3	6	2	9	0.8	99.90	99.70	100.10
50	50	2	4	4	2	0.1	66.10	63.16	69.04
50	50	2	4	4	2	0.4	97.90	97.01	98.79

Ν	TimeP	NFAC	NVAR	NFAC2	NVAR2	Autoreg	Accuracy	Lower 95% C.I	Upper 95% C.I
50	50	2	4	4	2	0.8	99.90	99.70	100.10
50	50	2	6	3	4	0.1	100.00	100.00	100.00
50	50	2	6	3	4	0.4	100.00	100.00	100.00
50	50	2	6	3	4	0.8	100.00	100.00	100.00
50	50	3	4	2	6	0.1	95.90	94.67	97.13
50	50	3	4	2	6	0.4	95.30	93.99	96.61
50	50	3	4	2	6	0.8	94.50	93.08	95.92
50	50	3	6	2	9	0.1	86.60	84.49	88.71
50	50	3	6	2	9	0.4	99.80	99.52	100.08
50	50	3	6	2	9	0.8	100.00	100.00	100.00
50	100	2	4	4	2	0.1	66.80	63.88	69.72
50	100	2	4	4	2	0.4	99.20	98.65	99.75
50	100	2	4	4	2	0.8	100.00	100.00	100.00
50	100	2	6	3	4	0.1	100.00	100.00	100.00
50	100	2	6	3	4	0.4	100.00	100.00	100.00
50	100	2	6	3	4	0.8	100.00	100.00	100.00
50	100	3	4	2	6	0.1	99.80	99.52	100.08
50	100	3	4	2	6	0.4	99.90	99.70	100.10
50	100	3	4	2	6	0.8	99.50	99.06	99.94
50	100	3	6	2	9	0.1	96.60	95.47	97.73
50	100	3	6	2	9	0.4	100.00	100.00	100.00
50	100	3	6	2	9	0.8	100.00	100.00	100.00
100	50	2	4	4	2	0.1	69.80	66.95	72.65
100	50	2	4	4	2	0.4	99.40	98.92	99.88
100	50	2	4	4	2	0.8	100.00	100.00	100.00
100	50	2	6	3	4	0.1	100.00	100.00	100.00
100	50	2	6	3	4	0.4	100.00	100.00	100.00

Table 1Accuracy per conditions tested (continued)

Ν	TimeP	NFAC	NVAR	NFAC2	NVAR2	Autoreg	Accuracy	Lower 95% C.I	Upper 95% C.I
100	50	2	6	3	4	0.8	100.00	100.00	100.00
100	50	3	4	2	6	0.1	100.00	100.00	100.00
100	50	3	4	2	6	0.4	99.90	99.70	100.10
100	50	3	4	2	6	0.8	100.00	100.00	100.00
100	50	3	6	2	9	0.1	95.80	94.55	97.05
100	50	3	6	2	9	0.4	100.00	100.00	100.00
100	50	3	6	2	9	0.8	100.00	100.00	100.00
100	100	2	4	4	2	0.1	72.75	69.98	75.51
100	100	2	4	4	2	0.4	99.90	99.70	100.10
100	100	2	4	4	2	0.8	100.00	100.00	100.00
100	100	2	6	3	4	0.1	100.00	100.00	100.00
100	100	2	6	3	4	0.4	100.00	100.00	100.00
100	100	2	6	3	4	0.8	100.00	100.00	100.00
100	100	3	4	2	6	0.1	100.00	100.00	100.00
100	100	3	4	2	6	0.4	100.00	100.00	100.00
100	100	3	4	2	6	0.8	100.00	100.00	100.00
100	100	3	6	2	9	0.1	99.60	99.21	99.99
100	100	3	6	2	9	0.4	100.00	100.00	100.00
100	100	3	6	2	9	0.8	100.00	100.00	100.00

Table 1Accuracy per conditions tested (continued)

Table 1 and 4 show the mean accuracy per condition tested and the 95% confidence interval of the mean. Mean accuracy is defined as the average (across 500 simulated datasets per condition) number of times the EII was larger for the *NotEq* condition than for the *Eq* condition (in a scale from 0 to 100%). The grand mean accuracy was very high (Mean = 94.07%, SD = 23.62). The accuracy increases with an increase in the sample size, number of variables per factor, and as a function of the autoregressive parameter. When the number of



Figure 4. Mean accuracy per condition tested

factors is two, the EII presented a moderate accuracy for an autoregressive parameter of .1, 428 with only four variables per factor. As the number of variables per factor increased to six, 429 the accuracy was perfect. This is likely due to the configuration of the second group of 430 individuals in the *NotEq* condition. When the number of factors is two in the first group 431 with four variables per factor, the second group had four factors, each with two variables 432 only (total number of variables = 8). Increasing the number of variables to six in the first 433 group (with two main factors) means that the second group could have three factors with 434 four variables or four factors with three variables each (total number of variables = 12). So, 435 the six variables condition had better identified factors than the four variables condition 436

when the number of factors equals two for the second group of the NotEq condition. For the
three factors condition, sample size was more impactful when the number of variables is set
to four, and the autoregressive parameter was more impactful in the conditions with six
variables per factor.

Figure 5 shows the mean EII value per sample size, number of factors, number of 441 variables per factor, number of time points, and autoregressive parameter. Overall, the mean 442 EII of the NotEq condition is much higher than the mean of the Eq condition, showing that 443 the metric can reliably differentiate when the individuals have a similar structure compared 444 to when individuals have a different structure. This is an important evidence of the 445 reliability of the EII to capture the amount of information lost by representing a set of 446 measures as a single population (or interindividual) structure instead of representing the 447 measures as multiple individual structures. 448

Figure 6 shows the distribution of the EII value per number of factors and number of variables per factor (of the first group of individuals). Finally, Figure 7 shows the distribution of the EII value per number of variables per factor in the first group (rows) and in the second group of individuals (columns).

453 A bottom-up approach for finding generalizations

If the system is nonergodic, what then? From Figure 3, it's reasonable to think that elements of the network structure estimated for each individual can be used to search for similarities that could help uncover sub-groups of people or generalizable characteristics across individuals. Said differently, uncovering sub-groups could reveal ergodic systems underlying the overall system. Since Campbell and Stanley (1963) raised concerns regarding external validity in psychological research, generalizability hasn't received nearly enough attention (Yarkoni, 2022).

⁴⁶¹ The emphasis on generalizability has both weaknesses and strengths in the directions



Figure 5. Mean ergodicity information index (and 95 percent confidence interval of the mean) per condition tested



Figure 6. Distribution of the mean ergodicity fit index per number of factors (rows) and number of variables per factor (columns)

in which it has helped push research design and data analysis. Among the negatives, an 462 emphasis on drawing large, representative samples of participants has, to our way of 463 thinking, been detrimental to the study of complex psychological processes. The primary 464 reason for this assertion is that to the extent that processes are constructs, they may well 465 manifest themselves somewhat differently from individual to individual (Molenaar & 466 Nesselroade, 2012) making it meaningless to build an accurate representation of the process 467 with observable measures. Here, heterogeneity is not helpful. If one is trying to determine 468 the average number of bedrooms in single family dwellings in the U.S., then a truly 469 representative sampling of single family dwellings is highly desirable. But if one is trying to 470 determine the nature of the onset and progression of depression, a representative sample that 471



Figure 7. Distribution of the mean ergodicity fit index per number of variables per factor of the first group (rows) and number of variables per factor in the second group (columns)

⁴⁷² includes a variety of paths of onset and progression is not helpful and may aggregate over
⁴⁷³ paths to a point where the representation does not match any person in the sample.

In arguing for a "bottom-up" approach to the matter of generalizability. Nesselroade 474 and Molenaar (2016) argued: "It seems far more appropriate to apply tools that emphasize 475 first understanding individuals well and then identifying similarities across persons, thus 476 accruing generalizability gradually than initially fitting models to heterogeneous samples in 477 order to claim generalizability. Large, diverse samples of individuals may put a gleam in a 478 demographer's eye but a wide array of observed differences, however universal the underlying 479 mechanisms may be, can blind the behavioral scientist to the identification of general 480 processes (p.15)." Following this line of reasoning, we propose an approach that takes into 481 consideration the spectral properties of the individual network structures in the search for 482 sub-groups or generalizable characteristics using an information-theoretic metric to compare 483 the similarity between two structures. 484

485 Spectral properties of a network provide insights into not only the topological features

(i.e., connectivity between nodes) but also the community structure (Chauhan, Girvan, & 486 Ott, 2009) and temporal dynamics (Almendral & Diaz-Guilera, 2007). Recent work has 487 taken advantage of these properties to determine whether individual networks in a multiplex 488 network can be aggregated into groups (De Domenico, Nicosia, Arenas, & Latora, 2015). De 489 Domenico et al. (2015) proposed a multiplex network reduction approach by computing Von 490 Neuman entropy of two networks and computing their Jensen-Shannon Distance (see 491 Golino, Moulder, et al., 2020 for other applications of Von Neumann entropy in community 492 detection). Von Neumann entropy of a network can be computed as follows: 493

$$h_A = -\mathrm{Tr}[\mathcal{L}_G \log_2 \mathcal{L}_G],$$

where $\mathcal{L}_G = c \times (D - A)$ is the combinatorial Laplacian rescaled by c or one over the sum of the weights in the network. D is a matrix with the strength of each node (i.e., sum of each node's connections) on its respective diagonal and A is the network. \mathcal{L}_G is a density matrix that is then used to compute Von Neuman entropy:

$$h_A = -\sum_{i=1}^N \lambda_i \log_2(\lambda_i),$$

where λ are the eigenvalues of \mathcal{L}_G . Using Von Neumann entropy of the network, we can compute the Jensen-Shannon Divergence between two networks, which is a symmetric measure of dissimilarity related to Kullback-Leibler Divergence (De Domenico et al., 2015):

$$\mathcal{D}_{JS}(\rho||\sigma) = h(\mu) - \frac{1}{2}[h(\rho) + h(\sigma)],$$

where ρ and σ are \mathcal{L}_G of each network being compared and $\mu = \frac{1}{2}(\rho + \sigma)$. Taking the square root of \mathcal{D}_{JS} produces a [0, 1] bound metric often referred to as Jensen-Shannon Distance. Subtracting 1 from the Jensen-Shannon Distance produces a similarity metric rather than distance metric.

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De Domenico et al. (2015) used an entropy-based quality function and Ward's method 505 for agglomerative hierarchical clustering using the Jensen-Shannon Distance to determine 506 whether individual layers (i.e., networks) in a multiplex network can be aggregated. Our 507 approach follows allow similar lines: We use complete-linkage agglomerative hierarchical 508 clustering on the Jensen-Shannon Distance. The trees are then cut through all possible cuts, 509 obtaining as many sets of clusters as there are networks. Afterward, a similarity matrix (1 -510 Jensen-Shannon Distance) is obtained and used to compute modularity. Modularity is an 511 objective function used in many community detection algorithms in network science 512 (**newman2006modularity?**). The modularity metric quantifies the extent to which 513 within-cluster similarity is maximized and between-cluster similarity is minimized. The 514 clusters that maximize modularity are taken to be the groups of similar networks. 515

Data from 30 individuals, with 100 measurement points each and 12 variables, were 516 simulated using the direct autoregressive factor score model (**DAFS**; Engle & Watson, 1981; 517 Nesselroade et al., 2002). The individuals differed only in the number of underlying factors 518 (2, 3, or 4), forming three groups. Using DynEGA to estimate the intraindividual networks 519 and computing the Jensen-Shannon Distance for every pair of networks, enabled the 520 identification of three clusters of people (bottom left of Figure 8). On the top left of Figure 521 8, the hierarchical clustering dendogram with three groups correctly identified. On the top 522 right of Figure 8, the three different groups' network and community structures are 523 represented separately. On the bottom right of Figure 8, the population structure estimated 524 using DynEGA is shown for each group of individuals. 525

526

Testing EII and Determining Groups in Applied Examples

527 Bootsrtap EII Test

The EII provides a relative metric for the information lost when representing the sample as a between-person, population structure relative to within-person, individual



Figure 8

structures. Determining whether the amount of information lost is substantial requires understanding the information loss relative to when within-person structures match the between-person structure. One approach would be to generate random multivariate normal samples from the population structure's covariance matrix with the same number of participants and their corresponding time series to create a sampling distribution of EII to compare with the empirical EII. This approach, however, is computationally expensive (e.g., 100 bootstraps \times 50 participants \times 20 variables \times 10 time points).

A simpler approach is to acquire a distribution of EII when individuals are known to 537 be similar to the population structure. To do so, we rewire the population network structure 538 to generate variants that are close but not identical to the population structure. This 539 approach takes the population network structure and completely rewires an edge, randomly 540 reassigning an edge connecting two nodes to two other nodes. To obtain structures with 541 different levels of deviations away from the population structure, we randomly sample a 542 proportion of edges to be rewired from a uniform distribution with a minimum of 0.20 (20%) 543 of edges) to 0.40 (40% of edges). In addition, we add random noise to all edges in the 544 network randomly sampling from a uniform distribution ranging from -0.10 to 0.10. This 545 process simulates small to moderate perturbations to the population structure to obtain 546 structures that might be observed by individuals within a population. 547

We generate as many population variants as their are participants in the empirical 548 sample. With the new sample containing the population network variants, we compute EII 549 with the original population network as the population network and the population network 550 variants as the individuals. We repeat this process for X iterations (e.g., 100). This 551 approach creates a sampling distribution of EII that would expected when the individuals in 552 the population have slight deviations on the population structure. When the empirical EII is 553 greater than 97.5% of the samples (two-tailed p = 0.05), then there is significant information 554 lost when representing the sample as an aggregate, population network. In all other cases, 555 including when the empirical EII is less than 97.5% of the sample, the ergodic property holds 556

⁵⁵⁷ such that non-substantial information is lost when representing the sample as an aggregate, ⁵⁵⁸ population network. Applying this approach to the data generated in Figure ??, the data ⁵⁵⁹ are nonergodic when combined into an aggregate population but ergodic when separated by ⁵⁶⁰ cluster.

561 Single Cluster Test

If the bootstrap EII test is significant, then the next step is to determine whether there 562 are clusters (or groupings) that can be identified in the data. In some cases, there may be 563 mixtures of network structures that form groups of people that have similar structures. In 564 other cases, each person may be unique and there are no clusters. In the latter case, a 565 challenge arises: the JSD values between each network structure will be relatively 566 equidistant. Because of the similar distances between networks, many clustering algorithms 567 will suggest that one cluster exists. If the bootstrap EII test is significant, this result is 568 paradoxical: The data do not possess the ergodic property yet they form a single cluster. 569 While the bootstrap EII could be leveraged to interpret the result as the single cluster 570 reflects the lack of clustering, the evidence is unsatisfying. 571

We derive an additional test to determine whether a single cluster represents a single 572 homogeneous cluster or if a single cluster represents heterogeneity such that each individual 573 represents their own cluster (i.e., singleton clusters). This additional procedure is only 574 applied when the number of clusters is determined to equal one cluster. The procedure 575 begins by generating a random network with the same number of nodes and edges as each 576 individual network in the empirical sample. Next, JSD is computed between every pair of 577 random networks (as is done to the empirical networks). The JSD values of the empirical 578 and random networks are then compared using a paired samples t-test, comparing the 579 corresponding JSD values between the corresponding empirical (e.g., empirical network j and 580 empirical network j) and random networks (e.g., random network i and random network j). 581

If the empirical network JSD values are, on average, greater than or equal to the random network JSD values, then the empirical network similarities are no different than the random network similarities suggesting that the single cluster is not meaningful and no meaningful clusters exist. If the empirical JSD values are, on average, significantly less than the random network JSD values (based on a large effect adjusted for sample size; Pérez & Pericchi, 2014), then the empirical network similarities are substantially less than the random networks similarities suggesting that the single cluster is meaningful.

We performed a small simulation to determine the effectiveness of this single cluster 589 test. One hundred heterogeneous samples of thirty cases with fifty time points were 590 generated using the **DAFS** method. Each case in each sample was randomly generated with 591 36 variables that loaded (between 0.70 and 1.00) onto 2, 3, 4, 6, or 9 factors with moderate 592 to large error (0.20-0.50). Similarly, one hundred homogeneous samples were equivalent 593 generated except that, rather than each case being generated by a different factor structure, 594 all cases for each sample were generated from the same factor structure with no to moderate 595 error (0.00-0.30). For the heterogeneous samples, an iteration was thrown out and replaced if 596 there were two or more clusters identified. Given the conditions randomly sample from only 597 five potential factor structures (with the same number of items per factor), then its possible 598 that these clusters truly exist in the data and therefore are not representative of a 599 heterogeneous sample. For the homogeneous samples, an iteration was thrown out and 600 replaced if the bootstrap EII test was significant. The accuracy for correctly identifying 601 when heterogeneous samples were heterogeneous (i.e., truly singleton clusters) was 100% and 602 the accuracy for correctly identifying when homogeneous samples were homogeneous (i.e., 603 truly one cluster) was 97%. 604

With this single cluster test, all evidence can be used together. If the bootstrap EII test is significant and the single cluster test suggests singleton clusters, then the data are not ergodic and they possess no meaningful groupings. If the bootstrap EII test is non-significant and the single cluster test suggests a single cluster, then the data are ergodic and represent a single, homogeneous sample. What if the bootstrap EII test is significant and the single cluster test suggests a single cluster? While we expect this last circumstance to be rare, it's possible. Our stance is that the default position should be to assume that the data do *not* possess the ergodic property—that is, we believe evidence must be accumulated for ergodicity and a homogeneous sample. In null hypothesis testing terms, the null hypothesis should be that the data are not ergodic and do not represent a single, homogeneous sample and the alternative hypothesis requiring significant evidence to the contrary.

616 Empirical Examples

To demonstrate the bootstrap EII and single cluster tests, we evaluated two empirical 617 data examples that are commonly represented with an aggregate structure: personality and 618 brain networks. For personality, we used an empirical example taken from an intensive 619 longitudinal experience sampling study examining Big Five personality measured by the Big 620 Five Inventory-2 (Beck & Jackson, 2022; Soto & John, 2017). There were 199 participants 621 who completed between 1 and 158 time points. To ensure optimal data quality, we only 622 included participants who completed at least 20 time points and had network densities of at 623 least 0.15 (i.e., at least 15% of all possible connections present). These criteria narrowed the 624 final sample to 122 participants. For the brain data, we used resting state neuroimaging data 625 taken from a study examining creativity (Beaty et al., 2018), which used the 268-node Shen 626 brain atlas (Shen, Tokoglu, Papademetris, & Constable, 2013). We analyzed the time series 627 data of 176 participants. The resting state scan was 5 minutes in duration resulting in 150 628 time points (300 seconds with TR of 2 seconds). 629

Starting with personality, the bootstrap EII test was significant (EII = 1.417, $p = 0.02, EII_{bootstrap} = [1.405, 1.408]$) suggesting that the BFI-2 data were nonergodic. Following up the bootstrap EII test, the information clustering was applied and the single cluster test was performed. The single cluster test suggested that the empirical networks had significantly greater JSD values than the random networks $(M_{difference} = 0.11, t(7380) = 66.85, p < .001, d = 1.08)$ meaning that the single cluster detected was not meaningful and each individual in the sample is unique.

Like the personality data, the brain networks had a significant bootstrap EII test ($EII = 6.237, p = 0.02, EII_{bootstrap} = [5.855, 5.943]$) and significantly larger JSD values than random networks, $M_{difference} = 0.02, t(15399) = 159.28, p < .001, d = 1.78$. These results, in line with the personality data, suggest that (resting state) brain networks are not ergodic and no meaningful groups can be formed (i.e., each individual is unique). In sum, the personality and (resting state) brain networks do not possess ergodicity and therefore lose information when data are aggregated into a single population.

644

Discussion

For well over a century the study of variation has been the backbone of psychologists' 645 efforts to understand behavior and behavior change. Whether created via experimental 646 manipulation or measured as it exists in nature (Cronbach, 1957; e.g., Cronbach, 1975), 647 variation is the ore that has been dug up, assayed, weighed, and otherwise analyzed by 648 behavioral prospectors hoping to strike it rich. Between-person or interindividual variation 649 has been and is strongly favored in psychological research. Often, researchers infer 650 within-person or idiographic processes based on between-person variation including many 651 prominent theories. The reliance on interindividual variation has led to a field that appears 652 to implicitly assume ergodicity is a fundamental property of psychological processes. Despite 653 this ubiquitous assumption, many researchers have chosen to focus on how the behavior of 654 an individual could vary from one measurement occasion to the next (Beck & Jackson, 2022; 655 Fisher et al., 2018). 656

Intraindividual variability is no longer studied by a curious few. It has grown into a prominent psychological method over the last few decades, to the point that it warrants the undivided attention of both methodologists and substantive theorists. This article takes

steps toward more rigorous testing about whether researchers can safely ignore the different 660 patterns of intraindividual variability in order to construct more general interindividual 661 representations. We leveraged tools from dynamical systems, network science, and 662 information theory to develop tests to the determine cost of aggregation. DynEGA paired 663 with GLLA allows researchers to study how variables change together over time and the 664 general structure of their relations (Golino et al., 2022). The EII provides researchers with a 665 relative metric and bootstrap test to determine how much information is lost when 666 aggregating individuals into a single population structure and whether that amount is 667 significant. If significant information is lost, then the information clustering method can be 668 applied and, if necessary, a single cluster test can evaluate whether meaningful clusters exist 669 in the data. 670

We propose that EII quantifies the extent to which super-weak ergodicity holds in the 671 system. Super-weak ergodicity suggests that the individual structures of a system should 672 reflect, within reasonable error, the aggregate structure of the system. This level of 673 ergodicity is a minimum requirement of a system to be represented as an aggregate. Systems 674 that do not possess this property should not be aggregated because significant information is 675 lost—such that a mere fraction of the system can be expected to reflect the aggregate 676 system. In its present state, psychological processes are unlikely to be ergodic (Molenaar, 677 2004). Therefore, the measurement of ergodicity must be pursued in perpetuity. Our 678 position, as well as the position of many others (e.g., Fisher et al., 2018; Molenaar, 2004), is 679 that psychological processes should be considered nonergodic until they are repeatedly 680 demonstrated otherwise. 681

Our empirical examples take two commonly aggregated psychological phenomena, personality and (resting state) brain activity, and examined the extent to which they lose information when aggregated into a single population network. We show, by no slim margin, that personality and brain networks are nonergodic. We further show that the distance between each individual's structure is more than would be expected if it were generated by an equivalent random structure. For personality, these result question the extent to which
the Big Five generalize to individuals (Borkenau & Ostendorf, 1998). For brain networks,
these results add to the growing body of evidence that dynamic brain networks reveal
greater idiosyncrasies between people (Hutchison et al., 2013; Lurie et al., 2020). Together,
these results suggest that each individual person shows distinct processes that are lost when
modeled as a single aggregate structure. In short, "no two people are alike" (@ Nesselroade
& Molenaar, 2016).

What are the consequences of our findings? In the best case, we show that the 694 personality and brain network samples examined in our study are nonergodic and are 695 perhaps due to sampling variability. There are many underlying factors that make this 696 interpretation plausible such as the specific personality scale used (BFI-2) or the brain 697 network task (or lack of task) or the demographics of the samples (i.e., predominantly college 698 students). Our findings, however, are not the first to find that individual people have resting 690 state brain signatures that make them uniquely identifiable (e.g., Finn et al., 2015). In the 700 worst case, we show that personality and brain activity are nonergodic and that research 701 examining their aggregates, the vast majority of their literatures, rest on a faulty assumption. 702 If the latter is true, then our findings challenge the validity of the conclusions drawn by 703 decades of interindividual research in personality and neuroscience. Our hope is that the 704 truth is somewhere in between. 705

Psychology aims to understand the thoughts, emotions, and behavior of the person and 706 people. Despite repeated calls and manifestos, the study of people continues to dominate 707 psychology. With modern technology, everyday thoughts, emotions, and behaviors of the 708 person have never been more accessible. As psychologists begin to emphasize intraindividual 709 processes over interindividual predictions, statistical models that answer questions about the 710 dynamics of systems are needed (Epskamp et al., 2018b; e.g., Gates & Molenaar, 2012; 711 Moulder, Martynova, & Boker, 2021; Sterba & Bauer, 2010). The present work provides one 712 statistical tool to understand whether intraindividual dynamics can be reasonably 713

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- ⁷¹⁵ sub-aggregations when they cannot. Together, these tools allow researchers to establish
- ⁷¹⁶ generalizability starting with the person rather than searching for it across people
- 717 (Nesselroade & Molenaar, 2016).

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