Interrogating theoretical models of neural computation with deep inference

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1 Abstract

A cornerstone of theoretical neuroscience is the circuit model: a system of equations that captures a hypothesized neural mechanism. Such models are valuable when they give rise to an experimentally observed phenomenon – whether behavioral or in terms of neural activity – and thus can offer insights into neural computation. The operation of these circuits, like all models, critically depends on the choices of model parameters. When analytic derivation of the relationship between model parameters and computational properties is intractable, approximate inference and simulation-based techniques are relied upon for scientific insight. We bring the use of deep generative models for probabilistic inference to bear on this problem, learning distributions of parameters that produce the specified properties of computation. By learning parameter distributions that produce computations – an emergent property, we introduce a novel methodology that is particularly well-suited to the stochastic dynamical systems models predominant in our field of theoretical neuroscience. We motivate this methodology with a worked example analyzing sensitivity in the stomatogastric ganglion. We then use it to reveal the key factors of variability in a model of primary visual cortex, gain a mechanistic understanding of rapid task switching in superior colliculus models, and scale inference of large low-rank RNN's exhibiting stable amplification. While much use of deep learning in theoretical neuroscience focuses on drawing analogies between optimized neural architectures and the brain, this work illustrates how we can further leverage the power of deep learning towards solving inverse problems in theoretical neuroscience.

2 Introduction

The fundamental practice of theoretical neuroscience is to use a mathematical model to understand neural computation, whether that computation enables perception, action, or some intermediate processing [1]. A neural computation is systematized with a set of equations – the model – and these equations are motivated by biophysics, neurophysiology, and other conceptual considerations. The function of this system is governed by the choice of model parameters, which when configured in a particular way, give rise to a measurable signature of a computation. The work of analyzing a model then requires solving the inverse problem: given a computation of interest, how can we reason about these particular parameter configurations? The inverse problem is crucial for reasoning about likely parameter values, uniquenesses and degeneracies, and predictions made by the model.

Consider the idealized practice: one carefully designs a model and analytically derives how model parameters govern the computation. Seminal examples of this gold standard (which often adopt approaches from statistical physics) include our field's understanding of memory capacity in associative neural networks [2], chaos and autocorrelation timescales in random neural networks [3], the paradoxical effect [4], and decision making [5]. Unfortunately, as circuit models include more biological realism, theory via analytical derivation becomes intractable. Alternatively, we can gain insight into these complex models by identifying all of the parameters consistent with the emergent phenomena of interest. By examining the structure of the full space of possible parameters, scientists can reason about the sensitivity and robustness of the model with respect to different parameter combinations [6, 7, 8, 9, 10].

The preferred formalism for parameter identification in science, statistical inference, has been used to great success in neuroscience through the stipulation of statistical generative models [11, 12, 13, 14, 15, 16, 17, 18, 19, 20, 21, 22, 23, 24, 25] (see review, [26]). However, most neural circuit models in theoretical neuroscience stipulate a noisy system of differential equations that can only be sampled or realized through forward simulation; they lack the explicit likelihood central to the probabilistic modeling toolkit. Therefore, the most popular approaches to the inverse problem have been likelihood-free methods such as approximate Bayesian computation (ABC) [27, 28], in which a set of reasonable parameters estimates is obtained via simulation and rejection.

Of course, the challenge of doing inference in complex models has arisen in many scientific fields. In response, the machine learning community has made remarkable progress in recent years, via the use of deep neural networks as a powerful inference engine: a flexible function family that can

map observations back to probability distributions quantifying the likely parameter configurations. One celebrated example of this approach from machine learning, of which we draw key inspiration for this work, is the variational autoencoder (VAE) [29, 30], which uses a deep neural network to induce an (approximate) posterior distribution on hidden variables in a latent variable model, given data. Indeed, these tools have been used to great success in neuroscience as well, in particular for interrogating parameters (sometimes treated as hidden states) in models of both cortical population activity [31, 32, 33, 34] and animal behavior [35, 36, 37]. These works have used deep neural networks to expand the expressivity and accuracy of statistical models of neural data [26].

Existing approaches to the inverse problem in theoretical neuroscience fall short in three key ways. First, theoretical models of neural computation aim to reflect a complex biological reality, and as a result, such models lack tractable likelihoods. Neuroscientists therefore resort to using approximate Bayesian computation, which requires a rejection heuristic, scales poorly, and only obtains sets of non-rejected parameters lacking probabilities. Second, is the undesirable trade-off between the flexibility and sampling speed of the approximated posterior distributions. Sampling-based approaches to statistical inference (e.g. ABC and Markov chain Monte Carlo (MCMC)) have flexibility in approximation, but must be executed continually for increasing samples. While variational approaches often result in fast sampling and sensitivity measurements post-optimization, existing approaches have relied on simplified classes of distributions. These simple distributions (e.g. mean-field gaussians) restrict the flexibility of the posterior approximation. And third, one can never assume what inferred model parameters may predict. This is well understood when considering Box's loop and the role of posterior predictive checks in the development and critique of scientific models [38, 39]. Uncertainty about the properties of inferred model predictions introduce a conceptual degree of freedom to the inverse problem that may be unnecessary and undesirable given the scientific motivation.

To address these three challenges, we developed an inference methodology – 'emergent property inference' – which learns a distribution over parameter configurations in a theoretical model. This distribution has two critical properties: *(i)* it is chosen such that draws from the distribution (parameter configurations) correspond to systems of equations that give rise to a specified emergent property (a set of constraints); and *(ii)* it is chosen to have maximum entropy given those constraints, such that we identify all likely parameters and can use the distribution to reason about parametric sensitivity and degeneracies [40]. First, we use stochastic gradient techniques in the spirit of likelihood-free variational inference [41] to enable inference in likelihood-free models of

neural computation. Second, we stipulate a bijective deep neural network that induces a flexible family of probability distributions over model parameterizations with a probability density we can calculate [42, **?**, 43], which confers fast sampling and sensitivity measurements. Third, we quantify the notion of emergent properties as a set of moment constraints on datasets generated by the model. Thus, an emergent property is not a single data realization, but a phenomenon or a feature of the model, which is ultimately the object of interest in theoretical neuroscience. Conditioning on an emergent property requires a variant of deep probabilistic inference methods, which we have previously introduced [44]. Taken together, emergent property inference (EPI) provides a methodology for inferring parameter configurations consistent with a particular emergent phenomena in theoretical models. We use a classic example of parametric degeneracy in a biological system, the stomatogastric ganglion [45], to motivate and clarify the technical details of EPI.

Equipped with this methodology, we then investigated three models of current importance in theoretical neuroscience. These models were chosen to demonstrate generality through ranges of biological realism (from conductance-based biophysics to recurrent neural networks), neural system function (from pattern generation to decision making), and network scale (from four to hundreds of neurons). First, we use EPI to understand the characteristics of noise that govern Fano factor in a stochastic four neuron-type model of primary visual cortex. Second, we discover connectivity patterns in superior colliculus resilient to optogenetic perturbation by using EPI to condition on rapid task switching. The novel scientific insights offered by EPI contextualize and clarify the previous studies exploring these models [46, 47]. Third, we emphasize the methodological advancement of EPI by inferring high-dimensional distributions of RNN connectivities exhibiting stable amplification. These results point to the value of deep inference for the interrogation of biologically relevant models.

3 Results

3.1 Motivating emergent property inference of theoretical models

Consideration of the typical workflow of theoretical modeling clarifies the need for emergent property inference. First, one designs or chooses an existing model that, it is hypothesized, captures the computation of interest. To ground this process in a well-known example, consider the stomatogastric ganglion (STG) of crustaceans, a small neural circuit which generates multiple rhythmic muscle activation patterns for digestion [48]. Despite full knowledge of STG connectivity and a

precise characterization of its rhythmic pattern generation, biophysical models of the STG have complicated relationships between circuit parameters and neural activity [45, 7]. A subcircuit model of the STG [49] is shown schematically in Figure 1A, and note that the behavior of this model will be critically dependent on its parameterization – the choices of conductance parameters $\mathbf{z} = [g_{\text{el}}, g_{\text{synA}}]$. Specifically, the two fast neurons ($f1$ and $f2$) mutually inhibit one another, and oscillate at a faster frequency than the mutually inhibiting slow neurons ($s1$ and $s2$). The hub neuron (hub) couples with either the fast or slow population or both.

Second, once the model is selected, one defines the emergent phenomena of scientific interest. In the STG example, we are concerned with neural spiking frequency, which emerges from the dynamics of the circuit model 1B. An interesting emergent property of this stochastic model is when the hub neuron fires at an intermediate frequency between the intrinsic spiking rates of the fast and slow populations. This emergent property is shown in Figure 1C at an average frequency of 0.55Hz.

Third, parameter analyses ensue: brute-force parameter sweeps, ABC sampling, and sensitivity analyses are all routinely used to reason about what parameter configurations lead to an emergent property. In this last step lies the opportunity for a precise quantification of the emergent property as a statistical feature of the model. Once we have such a methodology, we can infer a probability distribution over parameter configurations that produce this emergent property.

Before presenting technical details (in the following section), let us understand emergent property inference schematically: EPI (Fig. 1D) takes, as input, the model and the specified emergent property, and as its output, produces the parameter distribution EPI (Fig. 1E). This distribution – represented for clarity as samples from the distribution – is then a scientifically meaningful and mathematically tractable object. In the STG model, this distribution can be specifically queried to reveal the prototypical parameter configuration for network syncing (the mode; Figure 1E yellow star), and how network syncing decays based on changes away from the mode. The eigenvectors (of the Hessian of the distribution at the mode) quantitatively formalize the robustness of unified intermediacy (Fig. 1B solid (v_1) and dashed (v_2) black arrows). Indeed, samples equidistant from the mode along these EPI-identified dimensions of sensitivity (v_1) and degeneracy (v_2) agree with error contours (Fig. 1B contours) and have diminished or preserved network syncing, respectively (Fig. 1F activity traces, Fig. S TODO) (see Section 5.2.1).

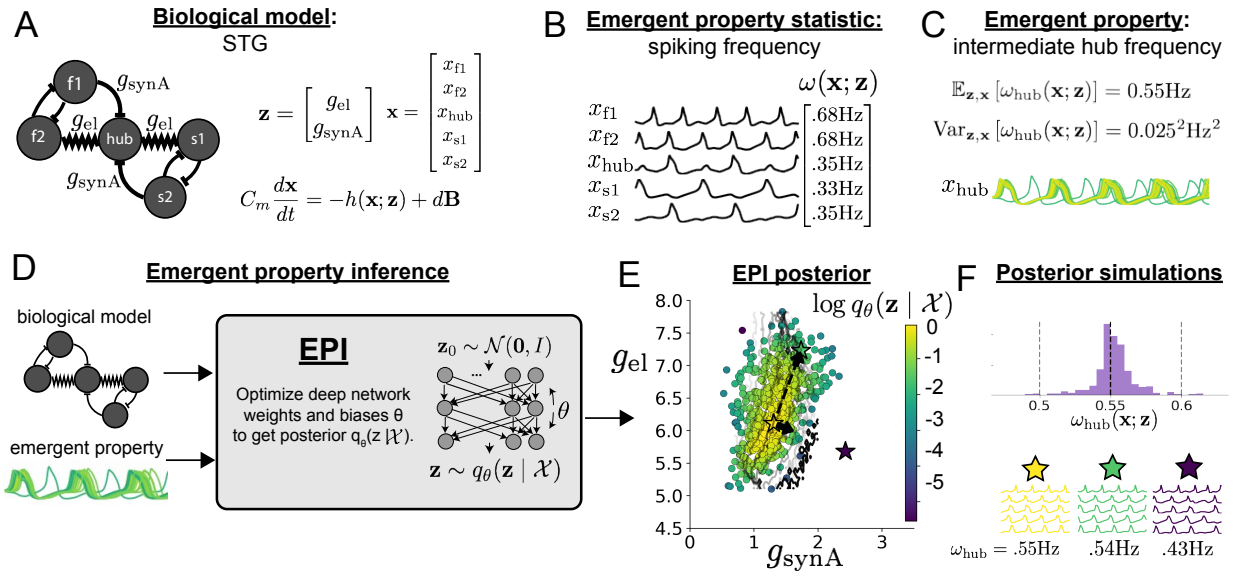

Figure 1: Emergent property inference (EPI) in the stomatogastric ganglion. **A**. Conductance-based biophysical model of the STG subcircuit. In the STG model, jagged connections indicate electrical coupling having electrical conductance g_{el}. Other connections in the diagram are inhibitory synaptic projections having strength g_{synA} onto the hub neuron, and $g_{\text{synB}} = 5\text{nS}$ for mutual inhibitory connections. Parameters are represented by the vector $\mathbf{z}$ and membrane potentials by the vector $\mathbf{x}$. The evolution of this model's activity $\mathbf{x}(t)$ is predicated by differential equations. **B**. Spiking frequency $\omega(\mathbf{x}; \mathbf{z})$ is an emergent property statistic. Spiking frequency is measured from simulated activity of the STG model at parameter choices of $g_{\text{el}} = 4.5\text{nS}$ and $g_{\text{synA}} = 3\text{nS}$. **C**. The emergent property of intermediate hub frequency, in which the hub neuron fires at a rate between the fast and slow frequencies. Simulated activity traces are colored by log probability density of their generating parameters in the EPI-inferred distribution (Panel E). **D**. For a choice of model and emergent property, emergent property inference (EPI) learns a distribution of the model parameters $\mathbf{z} = [g_{\text{el}}, g_{\text{synA}}]$ producing intermediate hub frequency. Deep probability distributions map a simple random variable $\mathbf{z}_0$ through a deep neural network with weights and biases $\boldsymbol{\theta}$ to parameters $\mathbf{z} = g_{\boldsymbol{\theta}}(\mathbf{z}_0)$ distributed as $q_{\boldsymbol{\theta}}(\mathbf{z} \mid \mathcal{X})$. In EPI optimization, stochastic gradient steps in $\boldsymbol{\theta}$ are taken such that entropy is maximized, and the emergent property $\mathcal{X}$ is produced. **E**. The EPI distribution of STG model parameters producing intermediate hub frequency. Samples are colored by log probability density. Distribution contours of hub neuron frequency from mean of .55 Hz are shown at levels of .525, .53, ... .575 Hz (dark to light gray away from mean). Frequencies are averages over the stochasticity of the model. Eigenvectors of the Hessian at the mode of the inferred distribution are indicated as $\mathbf{v}_1$ (solid) and $\mathbf{v}_2$ (dashed) with lengths scaled by the square root of the absolute value of their eigenvalues. Simulated activity is shown for three samples (stars). **F** Simulations from parameters in E. (Top) The predictive distribution of the posterior obeys the constraints stipulated by the emergent property. The black and gray dashed lines show the mean and two standard deviations according the emergent property, respectively. (Bottom) Simulations at the starred parameter values.

3.2 A deep generative modeling approach to emergent property inference

Emergent property inference (EPI) systematizes the three-step procedure of the previous section. First, we consider the model as a coupled set of differential equations [49]. In the running STG example, the model activity $\mathbf{x} = [x_{\text{f1}}, x_{\text{f2}}, x_{\text{hub}}, x_{\text{s1}}, x_{\text{s2}}]$ is the membrane potential for each neuron, which evolves according to the biophysical conductance-based equation:

$$C_m \frac{d\mathbf{x}(t)}{dt} = -h(\mathbf{x}(t); \mathbf{z}) + d\mathbf{B} \tag{1}$$

where C_m=1nF, and $\mathbf{h}$ is a sum of the leak, calcium, potassium, hyperpolarization, electrical, and synaptic currents, all of which have their own complicated dependence on $\mathbf{x}$ and $\mathbf{z} = [g_{\text{el}}, g_{\text{synA}}]$, and $d\mathbf{B}$ is white gaussian noise (see Section 5.2.1).

Second, we define the emergent property, which as above is "intermediate hub frequency" (Figure 1C). Quantifying this phenomenon is straightforward: we stipulate that the hub neuron's spiking frequency – denoted $\omega_{\text{hub}}(\mathbf{x})$ is close to an intermediate frequency of 0.55Hz. Mathematically, we achieve this via constraints on the mean and variance of the hub neuron spiking frequency.

$$\begin{aligned} \mathcal{X} \quad : \quad \mathbb{E}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x};\mathbf{z})\right] \quad &\triangleq \quad \mathbb{E}_{\mathbf{z},\mathbf{x}}\left[\omega_{\text{hub}}(\mathbf{x};\mathbf{z})\right] \quad = \quad \left[0.55\right] \quad \triangleq \quad \boldsymbol{\mu} \\ \text{Var}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x};\mathbf{z})\right] \quad &\triangleq \quad \text{Var}_{\mathbf{z},\mathbf{x}}\left[\omega_{\text{hub}}(\mathbf{x};\mathbf{z})\right] \quad = \quad \left[0.025^2\right] \quad \triangleq \quad \boldsymbol{\sigma}^2. \end{aligned} \tag{2}$$

The emergent property statistic $f(\mathbf{x};\mathbf{z}) = \omega_{\text{hub}}(\mathbf{x};\mathbf{z})$ along with its constrained mean $\boldsymbol{\mu}$ and variance $\boldsymbol{\sigma}^2$ define the emergent property denoted $\mathcal{X}$.

Third, we perform emergent property inference: we find a distribution over parameter configurations $\mathbf{z}$, and insist that samples from this distribution produce the emergent property; in other words, they obey the constraints introduced in Equation 2. This distribution will be chosen from a family of probability distributions $\mathcal{Q} = \{q_{\boldsymbol{\theta}}(\mathbf{z}) : \boldsymbol{\theta} \in \Theta\}$, defined by a deep generative distribution of the normalizing flow class [42, **?**, 43] – neural networks which transform a simple distribution into a suitably complicated distribution (as is needed here). This deep distribution is represented in Figure 1C (see Section 5.1). Then, mathematically, we must solve the following optimization program:

$$\begin{aligned} q_{\boldsymbol{\theta}}(\mathbf{z} \mid \mathcal{X}) \quad = \quad & \underset{q_{\boldsymbol{\theta}} \in \mathcal{Q}}{\text{argmax}}\, H(q_{\boldsymbol{\theta}}(\mathbf{z})) \\ & \text{s.t.} \quad \mathcal{X} : \quad \mathbb{E}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x};\mathbf{z})\right] = \boldsymbol{\mu}, \text{Var}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x};\mathbf{z})\right] = \boldsymbol{\sigma}^2 \end{aligned} \tag{3}$$

where $f(\mathbf{x}, \mathbf{z}), \boldsymbol{\mu}$, and $\boldsymbol{\sigma}$ are defined as in Equation 10. According to the emergent property of interest, $f(\mathbf{x}, \mathbf{z})$ may contain multiple statistics, in which case the mean and variance vectors $\boldsymbol{\mu}$ and $\boldsymbol{\sigma}^2$ match this dimension. Finally, we recognize that many distributions in $\mathcal{Q}$ will respect the emergent property constraints, so we select that which has maximum entropy. This principle, captured in Equation 3 by the primal objective H, identifies parameter distributions with minimal assumptions beyond some chosen structure [50, 51, 44, 52]. Such a normative principle of maximum entropy, which is also that of Bayesian inference, naturally fits with our scientific objective of reasoning about parametric sensitivity and robustness. The recovered distribution of EPI is as variable as possible along each parametric manifold such that it produces the emergent property.

EPI optimizes the weights and biases $\boldsymbol{\theta}$ of the deep neural network (which induces the probability distribution) by iteratively solving Equation 3. The optimization is complete when the sampled models with parameters $\mathbf{z} \sim q_{\boldsymbol{\theta}}(z \mid \mathcal{X})$ produce activity consistent with the specified emergent property (Fig. 8). Such convergence is evaluated with a hypothesis test that the means and variances of each emergent property statistic are not different than their constrained values (see Section 5.1.3). Further validation of EPI is available in the supplementary materials, where we analyze a simpler model for which ground-truth statements can be made (Section 5.1.4).

In relation to broader methodology, inspection of the EPI objective reveals a natural relationship to posterior inference. Specifically, EPI executes a novel variant of Bayesian inference with a uniform prior and a gaussian likelihood on the emergent property statistic (see Section 5.1.5). A key advantage of EPI over established Bayesian inference is that the predictions made by the inferred distribution are constrained to produce the specified emergent property. Equipped with this method, we may examine structure in posterior distributions or make comparisons between posteriors conditioned at different levels of the same emergent property statistic. In Sections 3.3 and 3.4, we prove out the value of EPI by using it to investigate and produce novel insights into two prominent models in neuroscience. Subsequently in Section 3.5, we show EPI's superiority in parameter scalability and fidelity of the posterior predictive distribution by conditioning on stable amplification in low-rank RNNs.

3.3 EPI reveals how noise across neural population types governs Fano factor in a stochastic inhibition stabilized network

Dynamical models of excitatory (E) and inhibitory (I) populations with supralinear input-output function have succeeded in explaining a host of experimentally documented phenomena. In a regime

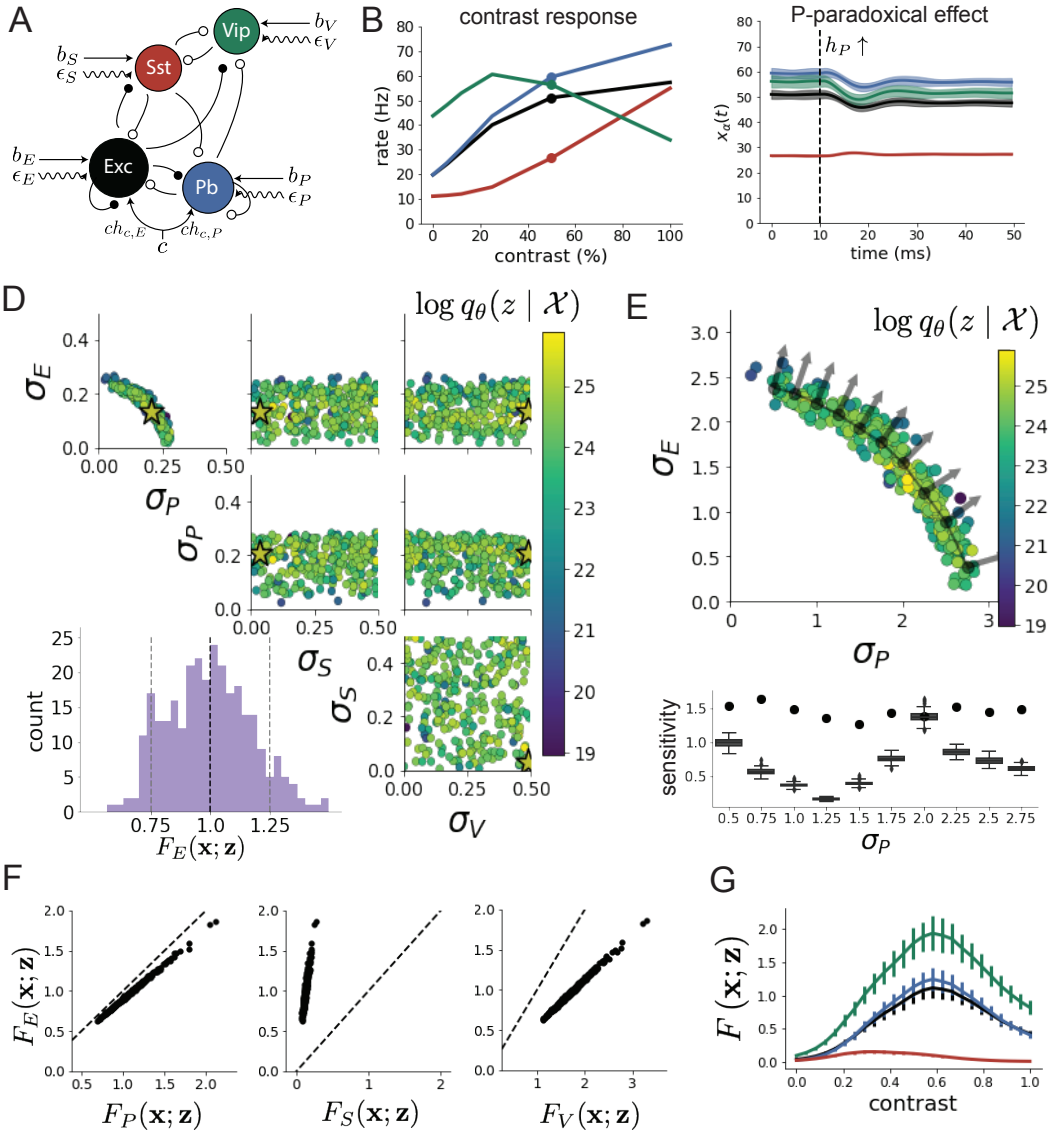

Figure 2: Emergent property inference of a stochastic stabilized supralinear network. **A**. Four-population model of primary visual cortex with excitatory (black), parvalbumin (blue), somatostatin (red), and VIP (green) neurons (excitatory and inhibitory projections filled and unfilled, respectively). Some neuron-types largely do not form synaptic projections to others ($|W_{\alpha_1,\alpha_2}| < 0.025$). Each neural population receives a baseline input $\mathbf{h}_b$, and the E- and P-populations also receive a contrast-dependent input $\mathbf{h}_b$. Additionally, each neural population receives a slow noisy input $\boldsymbol{\epsilon}$. **B**. Responses of the deterministic model ($\boldsymbol{\epsilon} = \mathbf{0}$) to varying contrasts. The response at 50% contrast (dots) is the focus of our analysis. **C**. Paradoxical response of the stochastic model to a small increase in input to the P-population. **D**. EPI posterior of noise parameters $\mathbf{z}$ conditioned on realistic E-population Fano factors. The posterior predictive distribution is show on the bottom-left. and the mode of the distribution is starred. **E**. (Top) Enlarged visualization of the σ_E-σ_P marginal distribution of the posterior. Each gray dot is a choice of σ_P, for which a constrained mode $z^*(\sigma_P, P)$ is chosen. The arrows show the most sensitive dimensions of the Hessian evaluated at these modes. (Bottom) Such sensitive dimensions of the Hessian (dots) are significantly more sensitive than randomly chosen dimensions (box and whiskers). **F**. The Fano factor of the E-population is strongly correlated with each other neuron-type population. **G**. Mean and standard deviation (across EPI posterior) of Fano factor of each neuron-type population at each level of contrast.

characterized by inhibitory stabilization of strong recurrent excitation, these models give rise to paradoxical responses [4], selective amplification [53, 54], surround suppression [55] and normalization [56]. Despite their strong predictive power, E-I circuit models rely on the assumption that inhibition can be studied as an indivisible unit. However, experimental evidence shows that inhibition is composed of distinct elements – parvalbumin (P), somatostatin (S), VIP (V) – composing 80% of GABAergic interneurons in V1 [57, 58, 59], and that these inhibitory cell types follow specific connectivity patterns (Fig. 2A) [60]. Recent theoretical advances [46, 61, 62], have only started to address the consequences of this multiplicity in the dynamics of V1, strongly relying on linear theoretical tools. Here, we use EPI to characterize the properties of slow noise in a stochastic version of this model, which result in biologically realistic responses.

We considered the contrast response of a nonlinear dynamical V1 circuit model (Fig. 2A) with a state comprised of each neuron-type population's rate $\mathbf{x} = [x_E, x_P, x_S, x_V]^\top$. Each population receives recurrent input $W\mathbf{x}$ from synaptic projections of effective connectivity W and an external input $\mathbf{h}$, which determine the population rate via nonlinearity $\phi = []_+^2$ (see Section 5.2.2). The circuit model evolves form an initial condition $\mathbf{x}(0) \sim \mathcal{U}([10, 25])$ with time constant $\tau = 1$ms according to a contrast-dependent input $\mathbf{h}$ and slow noise $\boldsymbol{\epsilon}$ of time constant $\tau_{\text{noise}} = 5$ms. This model is the stochastic stabilized supralinear network (SSSN) [63] generalized to have inhibitory multiplicity

$$\tau \frac{d\mathbf{x}}{dt} = -\mathbf{x} + \phi(W\mathbf{x} + \mathbf{h} + \boldsymbol{\epsilon}). \tag{4}$$

As contrast increases, input to the E- and P-populations increases relative to a baseline input $\mathbf{h}_b$ via $\mathbf{h}_c$

$$\mathbf{h} = \mathbf{h}_b + c\mathbf{h}_c, \tag{5}$$

where $h_{c,E}, h_{c,P} > 0$ and $h_{c,S}, h_{c,V} = 0$. In this analysis, we fixed W,$\mathbf{h}_b$, and $\mathbf{h}_c$ to values fit to mean contrast responses in mice with the deterministic model [64] ($\boldsymbol{\epsilon} = \mathbf{0}$, Fig. 2B, see Section 5.2.2). At all contrasts, the E-population of this SSSN is unstable without recurrent inhibitory feedback. At 50% contrast, only the P-population exhibits the paradoxical effect (2C, Fig. 9), so the network is P-stabilized.

The slow noise of the SSSN is an Ornstein-Uhlenbeck process

$$\tau_{\text{noise}} d\epsilon_\alpha = -\epsilon_\alpha dt + \sqrt{2\tau_{\text{noise}}}\sigma_\alpha dB, \tag{6}$$

parameterized by σ_α, which can be different for each neuron type,

$$\mathbf{z} = [\sigma_E, \sigma_P, \sigma_S, \sigma_V]^\top. \tag{7}$$

For this SSSN, we are interested in the parameters of slow noise that produce realistic stochastic fluctuations. Here, we quantify this emergent property as having an excitatory population Fano factor near 1:

$$\begin{aligned} \mathcal{X} \quad : \quad \mathbb{E}_{\mathbf{z}}\Big[F_E(\mathbf{x};\mathbf{z})\Big] &= 1 \\ \mathrm{Var}_{\mathbf{z}}\Big[F_E(\mathbf{x};\mathbf{z})\Big] &= 0.125^2, \end{aligned} \tag{8}$$

where $F_\alpha(\mathbf{x};\mathbf{z})$ is the Fano factor of the α-population.

We ran EPI to obtain a posterior $q_{\boldsymbol{\theta}}(\mathbf{z} \mid \mathcal{X})$, where each parameter $\mathbf{z}$ produces biologically realistic levels of E-population variability (Fig. 2D). From the marginal distribution of σ_E and σ_P (Fig. 2D, top-left), we can see that $F_E(\mathbf{x};\mathbf{z})$ is sensitive to the combination of σ_E and σ_P. In fact, the posterior obtained through EPI offers exactly how this sensitivity changes along this ridge of the posterior (Fig. 2E). σ_S and σ_V are degenerate with respect to $F_E(\mathbf{x};\mathbf{z})$ evidenced by the uniform distribution in those dimensions of the posterior (Fig. 2D, bottom-right). Together, this posterior indicates a parametric manifold of degeneracy with respect to Fano factor: the ridge visualized in the σ_E-σ_P marginal (Fig. 10) and the dimensions of σ_S and σ_V.

Greater σ_E and σ_P confer greater Fano factor, and the Fano factors of each neuron-type are strongly correlated across the posterior (Fig 2F), showing that Fano factor of each neuron-type can be modulated globally via σ_E and σ_P. Furthermore, across the entire posterior distribution of noise parameterizations, we find that when contrast is increased above 50%, variability is quenched for all neuron types (Fig 2G). In summary, we used EPI to obtain a posterior of SSSNs producing realistic Fano factors, which allowed degenerate manifold identification via sample visualization, fast sensitivity measurements via Hessian evaluation, and predictions of variability quenching.

3.4 EPI identifies neural mechanisms of flexible task switching

In a rapid task switching experiment [65], rats were explicitly cued on each trial to either orient towards a visual stimulus in the Pro (P) task or orient away from a visual stimulus in the Anti (A) task (Fig. 3A). Neural recordings in the midbrain superior colliculus (SC) exhibited two populations of neurons that simultaneously represented both task context (Pro or Anti) and motor response (contralateral or ipsilateral to the recorded side): the Pro/Contra and Anti/Ipsi neurons [47]. Duan et al. proposed a model of SC that, like the V1 model analyzed in the previous section, is a four-population dynamical system. We analyzed this model, where the neuron-type populations are functionally-defined as the Pro- and Anti-populations in each hemisphere (left (L) and right

(R)), their connectivity is parameterized geometrically (Fig. 3B). The input-output function of this model is chosen such that the population responses $\mathbf{x} = [x_{LP}, x_{LA}, x_{RP}, x_{RA}]^\top$ are bounded from 0 to 1 as a function ϕ of a dynamically evolving internal variable $\mathbf{u}$. The model responds to the side with greater Pro neuron activation; e.g. the response is left if $x_{LP} > x_{RP}$ at the end of the trial. The dynamics evolve with timescale $\tau = 0.09$ governed by connectivity weights W

$$\begin{aligned} \tau \frac{d\mathbf{u}}{dt} &= -\mathbf{u} + W\mathbf{x} + \mathbf{h} + d\mathbf{B} \\ \mathbf{x} &= \phi(\mathbf{u}) \end{aligned} \tag{9}$$

with white noise of variance 0.2^2. The input $\mathbf{h}$ is comprised of a cue-dependent input to the Pro or Anti populations, a stimulus orientation input to either the Left or Right populations, and a choice-period input to the entire network (see Section 5.2.3). Here, we use EPI to determine the changes in network connectivity $\mathbf{z} = [sW, vW, dW, hW]^\top$ resulting in execution of rapid task switching behavior.

We define rapid task switching behavior as accurate execution of each task. Inferred models should not exhibit fully random responses (50%), or perfect performance (100%), since perfection is never attained by even the best trained rats. We formulate rapid task switching as an emergent property by stipulating that the average accuracy in the Pro task $p_P(\mathbf{x}, \mathbf{z})$ and Anti task $p_A(\mathbf{x}, \mathbf{z})$ be 75% with variance $5\%^2$.

$$\begin{aligned} \mathcal{X} \quad : \quad \mathbb{E}_{\mathbf{z}} \begin{bmatrix} p_P(\mathbf{x};\mathbf{z}) \\ p_A(\mathbf{x};\mathbf{z}) \end{bmatrix} &= \begin{bmatrix} 75\% \\ 75\% \end{bmatrix} \\ \mathrm{Var}_{\mathbf{z}} \begin{bmatrix} p_P(\mathbf{x};\mathbf{z}) \\ p_A(\mathbf{x};\mathbf{z}) \end{bmatrix} &= \begin{bmatrix} 5\%^2 \\ 5\%^2 \end{bmatrix} \end{aligned} \tag{10}$$

A variance of $5\%^2$ performance in each task will confer a posterior producing performances ranging from about $65\% - 85\%$, allowing us to examine the properties of connectivity that yield better performance.

We ran EPI to obtain SC model connectivity parameters $\mathbf{z}$ producing rapid task switching (Fig. 3C). Some parameters were predictive of accuracy while others were not (Fig. 11), and often had different effects on p_P and p_A. To make sense of this inferred distribution, we took the eigendecomposition of the symmetric connectivity matrices $W = V\Lambda V^{-1}$, which results in the same basis vectors $\mathbf{v}_i$ for all W parameterized by $\mathbf{z}$ (Fig. 12A). These basis vectors have intuitive roles in processing for this task, and are accordingly named the *all* mode - all neurons co-fluctuate, *side* mode - one side dominates the other, *task* mode - the Pro or Anti populations dominate the

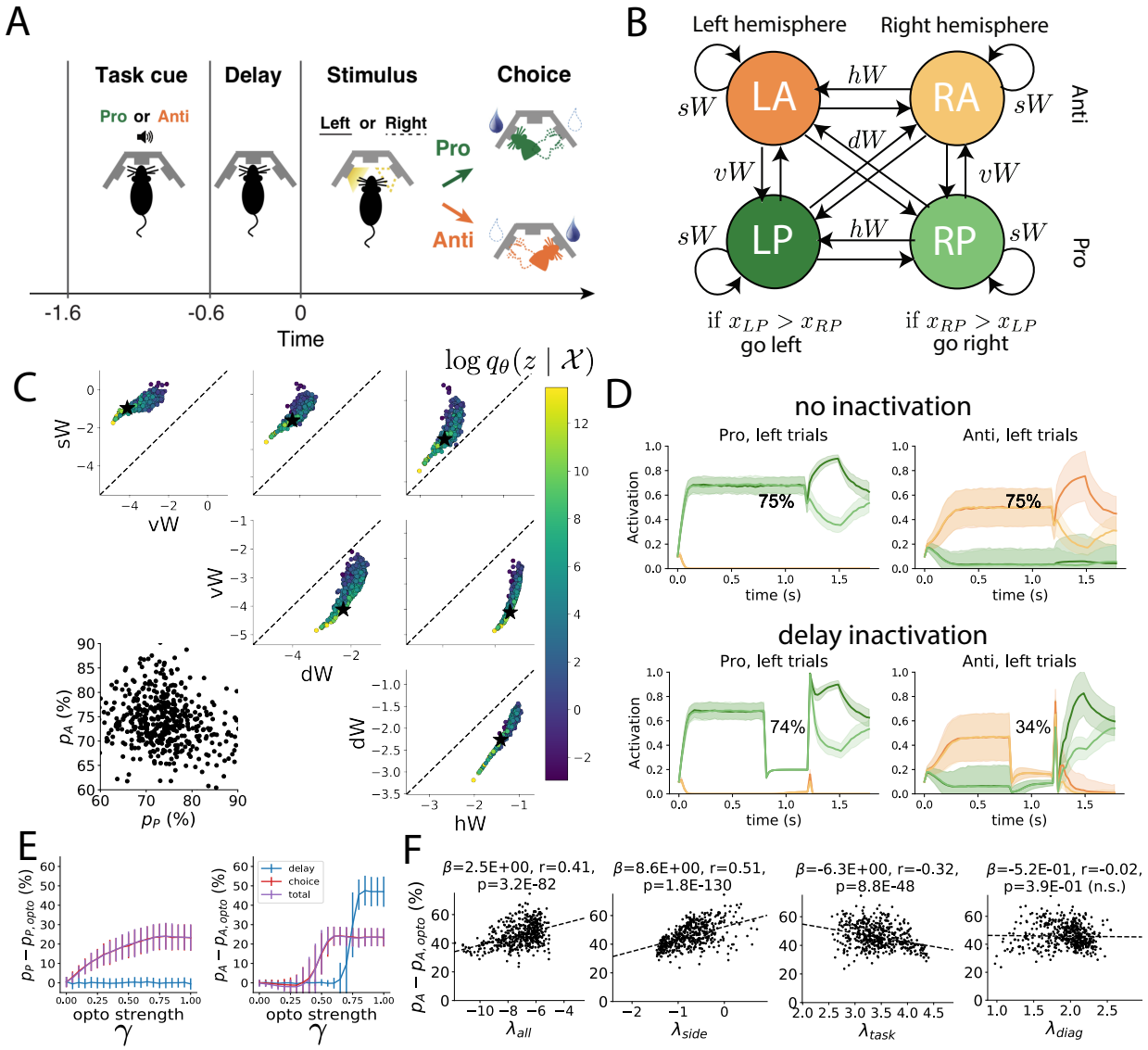

Figure 3: **A**. Rapid task switching behavioral paradigm (see text). **B**. Model of superior colliculus (SC). Neurons: LP - left pro, RP - right pro, LA - left anti, RA - right anti. Parameters: sW - self, hW - horizontal, vW -vertical, dW - diagonal weights. Subscripts P and A of connectivity weights indicate Pro or Anti populations. **C**. The EPI parameter distribution of rapid task switching networks. Black star indicates parameter choice of simulations (D). **D**. Simulations of an SC network from the EPI distribution with 75% accuracy in each task. Top row shows no inactivation during Pro and Anti trials, and bottom row shows simulations with delay period inactivation (optogenetic strength $\gamma = 0.7$). Shading indicates standard deviation across trials. **E**. Difference in performance of each task during inactivation. Inactivation level γ scales from no inactivation (0) to full inactivation (1). We compare delay period inactivation $1.2 < t < 1.5$ (blue), choice period inactivation $1.5 < t$ (red), and total inactivation $0 \leq t \leq 1.8$ (purple). **F**. The effect of delay period inactivation on Anti accuracy versus dynamics eigenvalues.

other, and *diag* mode - Pro- and Anti-populations of opposite hemispheres dominate the opposite pair.

Greater λ_{task}, λ_{side}, and λ_{diag} all produce greater Pro accuracy. This shows that strong task representations and hemispheric

dominance in the dynamics result in better execution of the Pro task. By visualizing these four variables together by p_A (Fig. 13B), we see that low λ_{task} and λ_{diag} producing strong Anti accuracy also have high λ_{side} and λ_{all}. Thus, stronger hemispheric dominance, relaxed task and diag mode dynamics, and slower circuit-wide decay result in greater Anti accuracy.

In agreement with experimental results from Duan et al., we found that inactivation above nominal strength during the delay period consistently decreased performance in the Anti task, but had no consistent effect on the Pro task (Fig. 3E) e.g. (Fig. 3D, bottom). This difference in resiliency across tasks to delay perturbation is a prediction made by the inferred EPI distribution, rather than an emergent property that was conditioned upon. Even though p_P and p_A are anticorrelated in the EPI posterior ($r = -0.15$, $p = 3.68 \times 10^{-}12$), greater p_P and p_A both result in decreased resiliency to delay perturbation in the Anti task (Fig. 14). Ultimately, lower λ_{side} and λ_{all} and greater λ_{task} produce networks more robust to delay perturbation in the Anti task (Fig. 3F)).

In summary, we used EPI to obtain the full distribution of connectivities that execute rapid task switching. This posterior revealed the mechanisms leading to greater accuracy in each task as well as those increasing resiliency to perturbation in the Anti task. Importantly, every connectivity from this inferred distribution predicts fragility and robustness of performance in the Anti and Pro tasks, respectively. EPI allows us to conclude that since *all* parameters of this model producing rapid task switching make such an experimentally verified prediction, we have a well-chosen model.

3.5 EPI scales well to high-dimensional parameter spaces

Here, we are interested in the scalability of EPI in number of parameters ($|\mathbf{z}|$). We consider rank-2 RNN with N neurons of connectivity

$$W = UV^\top + g\chi \tag{11}$$

and dynamics

$$\tau \frac{d\mathbf{x}}{dt} = -\mathbf{x} + W\mathbf{x} \tag{12}$$

where $U = \begin{bmatrix} \mathbf{u}_1 & \mathbf{u}_2 \end{bmatrix}$, $V = \begin{bmatrix} \mathbf{v}_1 & \mathbf{v}_2 \end{bmatrix}$, $\mathbf{u}_1, \mathbf{u}_2, \mathbf{v}_1, \mathbf{v}_2 \in [-1, 1]^N$, and $g = 0.01$.

We want to learn distributions of connectivity that produce stable amplification. Two conditions are both necessary and sufficient for RNNs to exhibit stable amplification [**?**]. These conditions are inequalities on real(λ_1) and λ_1^s the maximal real eigenvalue of W and the maximum eigenvalue of $W^s = \frac{W+W^\top}{2}$, respectively.

In our analysis, we seek to condition rank-2 networks of increasing size on a regime of stable amplification. Networks with real(λ_1) $= 0.5 \pm 0.5$ and $\lambda_1^s = 1.5 \pm 0.5$ will yield moderate amplification. EPI can naturally condition on this emergent property

$$\begin{aligned} \mathcal{X} \quad : \quad \mathbb{E}_{\mathbf{z},\mathbf{x}} \begin{bmatrix} \text{real}(\lambda_1) \\ \lambda_1^s \end{bmatrix} &= \begin{bmatrix} 0.5 \\ 1.5 \end{bmatrix} \\ \text{Var}_{\mathbf{z},\mathbf{x}} \begin{bmatrix} \text{real}(\lambda_1) \\ \lambda_1^s \end{bmatrix} &= \begin{bmatrix} 0.25^2 \\ 0.25^2 \end{bmatrix}. \end{aligned} \tag{13}$$

In contrast, SNPE cannot condition on the variance of observations across posterior. ' Thus, we condition on an observation x_0 located at the mean of our desired emergent property.

$$x_0 = \begin{bmatrix} \text{real}(\lambda_1) \\ \lambda_1^s \end{bmatrix} = \begin{bmatrix} 0.5 \\ 1.5 \end{bmatrix} \tag{14}$$

ABC methods define tolerance ϵ and distance for observed data x_0. Here, we chose $\epsilon = 0.5$, an $l-2$ distance, and the same choice for x_0 as in Equation 14.

EPI is capable of scaling to higher dimensional parameter spaces than ABC and SNPE. EPI consistently produces the same posterior predictive distribution independent of the dimensionality. SMC produces a limited variety of parameters due to the nature of its proposal generation algorithm, yet all parameters obtained produce stable amplification. SNPE's posterior predictive distribution is not necessarily close to the conditioning point, and is very dependent on dimensionality.

4 Discussion

NOTE: **This is the old discussion section. I will rewrite this based on our discussion of the rest of the draft.**

In neuroscience, machine learning has primarily been used to reveal structure in neural datasets [11, 12, 13, 14, 16, 18, 20, 22, 23, 24, 25] (see review, [26]). Such careful inference procedures

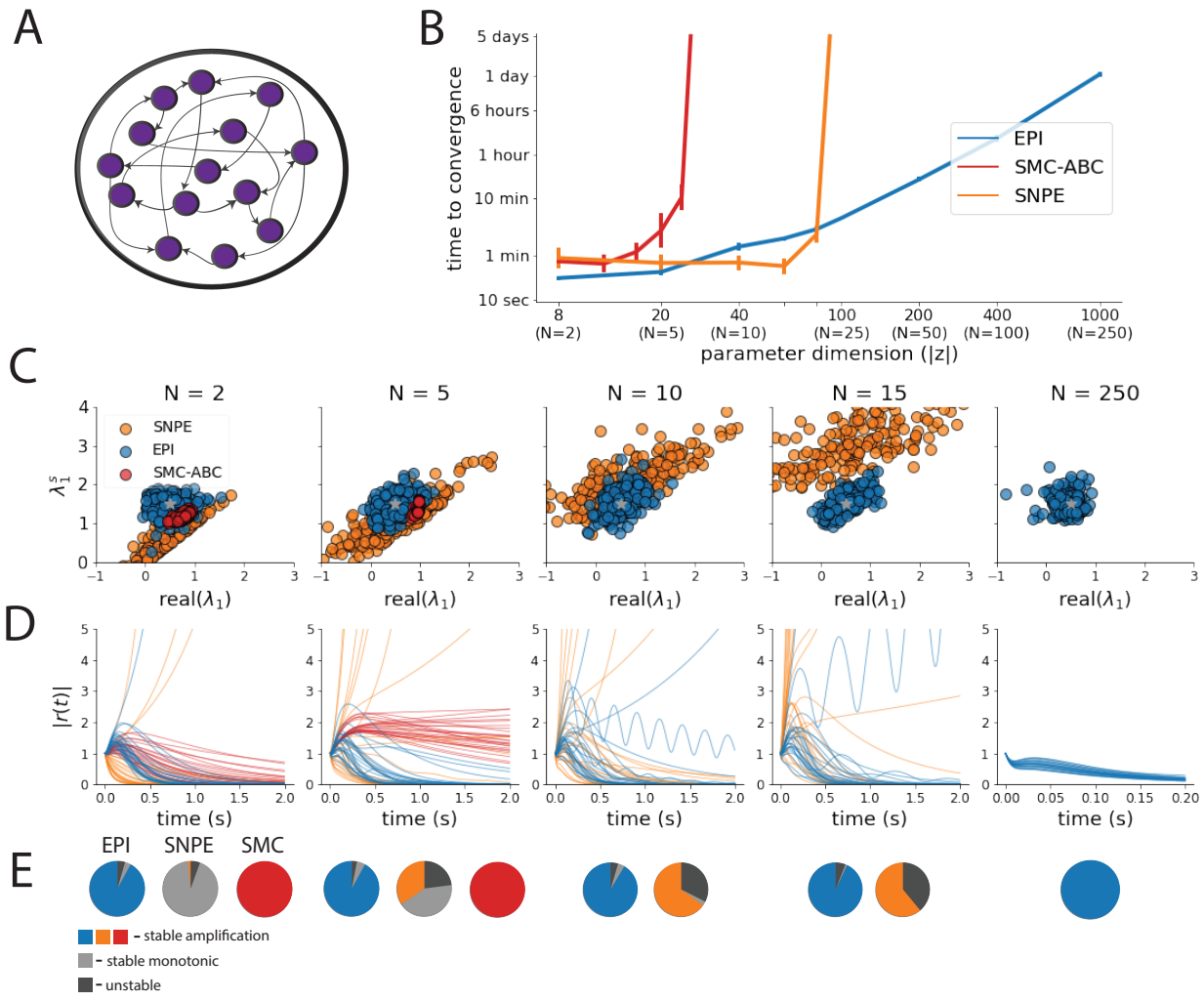

Figure 4: **A**. Recurrent neural network. **B**. EPI scales with z to high dimensions. Convergence definitions: EPI (blue) - satisfies all moment constraints, SNPE (orange)- produces at least $2/n_{\text{train}}$ parameter samples are in the bounds of emergent property (mean +- 0.5), and SMC-ABC (red) - 100 particles with $\epsilon <$ 0.5 are produced. **C**. Posterior predictive distributions of EPI (blue), SNPE (orange), and SMC-ABC (red). Gray star indicates emergent property mean, and gray dashed lines indicate two standard deviations corresponding to the variance constraint. For $N <= 6$ where SMC-ABC converges, samples are not diverse (path degeneracies). For $N >= 25$, SNPE does not produce a posterior approximation yielding parameters with simulations near x_0. **D**. Simulations of network parameters resulting from each method ($\tau = 100ms$). Each trace corresponds to simulation of one z. **E**. Ratio of obtained samples producing stable amplification.

are developed for these statistical models allowing precise, quantitative reasoning, which clarifies the way data informs beliefs about the model parameters. However, these statistical models lack resemblance to the underlying biology, making it unclear how to go from the structure revealed by these methods, to the neural mechanisms giving rise to it. In contrast, theoretical neuroscience has focused on careful mechanistic modeling and the production of emergent properties of computation. The careful steps of *i.*) model design and *ii.*) emergent property definition, are followed by *iii.*) practical inference methods resulting in an opaque characterization of the way model parameters govern computation. In this work, we replaced this opaque procedure of parameter identification in theoretical neuroscience with emergent property inference, opening the door to careful inference in careful models of neural computation.

Biologically realistic models of neural circuits often prove formidable to analyze. Two main factors contribute to the difficulty of this endeavor. First, in most neural circuit models, the number of parameters scales quadratically with the number of neurons, limiting analysis of its parameter space. Second, even in low dimensional circuits, the structure of the parametric regimes governing emergent properties is intricate. For example, these circuit models can support more than one steady state [66] and non-trivial dynamics on strange attractors [67].

In Section 3.3, we advanced the tractability of low-dimensional neural circuit models by showing that EPI offers insights about cell-type specific input-responsivity that cannot be afforded through the available linear analytical methods [46, 61, 62]. By flexibly conditioning this V1 model on different emergent properties, we performed an exploratory analysis of a *model* rather than a dataset, generating a set of testable hypotheses, which were proved out. Furthermore, exploratory analyses can be directed towards formulating hypotheses of a specific form. For example, model parameter dependencies on behavioral performance can be assessed by using EPI to condition on various levels of task accuracy (See Section 3.4). This analysis identified experimentally testable predictions (proved out *in-silico*) of patterns of effective connectivity in SC that should be correlated with increased performance.

In our final analysis, we presented a novel procedure for doing statistical inference on interpretable parameterizations of RNNs executing simple tasks. Specifically, we analyzed RNNs solving a posterior conditioning problem in the spirit of [68, 69]. This methodology relies on recently extended theory of responses in random neural networks with low-rank structure [70]. While we focused on rank-1 RNNs, which were sufficient for solving this task, this inference procedure generalizes to RNNs of greater rank necessary for more complex tasks. The ability to apply the probabilistic

model selection toolkit to RNNs should prove invaluable as their use in neuroscience increases.

EPI leverages deep learning technology for neuroscientific inquiry in a categorically different way than approaches focused on training neural networks to execute behavioral tasks [71]. These works focus on examining optimized deep neural networks while considering the objective function, learning rule, and architecture used. This endeavor efficiently obtains sets of parameters that can be reasoned about with respect to such considerations, but lacks the careful probabilistic treatment of parameter inference in EPI. These approaches can be used complementarily to enhance the practice of theoretical neuroscience.

TODO *merge this point in*

While much research in computational neuroscience has focused on optimizing neural architectures to process information and accomplish tasks [71], structure in parameter space of the set of optimized solutions is rarely discussed and lacks a probabilistic treatment. Talk about Wyktor's work here [72].

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Data availability statement:
The datasets generated during and/or analyzed during the current study are available from the corresponding author upon reasonable request.

Code availability statement:
The software written for the current study is available from the corresponding author upon reasonable request.

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5 Methods

5.1 Emergent property inference (EPI)

Determining the combinations of model parameters that can produce observed data or a desired output is a key part of scientific practice. Solving inverse problems is especially important in neuroscience, since we require complex models to describe the complex phenomena of neural computations. While much machine learning research has focused on how to find latent structure in large-scale neural datasets, less has focused on inverting theoretical circuit models conditioned upon the emergent phenomena they produce. Here, we introduce a novel method for statistical inference, which finds distributions of parameter solutions that only produce the desired emergent property. This method seamlessly handles neural circuit models with stochastic nonlinear dynamical generative processes, which are predominant in theoretical neuroscience.

Consider model parameterization $\mathbf{z}$, which is a collection of scientifically interesting variables that govern the complex simulation of data $\mathbf{x}$. For example (see Section 3.1), $\mathbf{z}$ may be the electrical conductance parameters of an STG subcircuit, and $\mathbf{x}$ the evolving membrane potentials of the five neurons. In terms of statistical modeling, this circuit model has an intractable likelihood $p(\mathbf{x} \mid \mathbf{z})$, which is predicated by the stochastic differential equations that define the model. Even so, we do not scientifically reason about how $\mathbf{z}$ governs all of $\mathbf{x}$, but rather specific phenomena that are a function of the data $f(\mathbf{x}; \mathbf{z})$. In the STG example, $f(\mathbf{x}; \mathbf{z})$ measures hub neuron frequency from the evolution of $\mathbf{x}$ governed by $\mathbf{z}$. With EPI, we learn distributions of $\mathbf{z}$ that results in an average and variance of $f(\mathbf{x}; \mathbf{z})$, denoted $\boldsymbol{\mu}$ and $\boldsymbol{\sigma}^2$. We refer to the collection of these statistical moments as an emergent property. Such emergent properties $\mathcal{X}$ are defined through choice of $f(\mathbf{x}; \mathbf{z})$ (which may be one or multiple statistics), $\boldsymbol{\mu}$, and $\boldsymbol{\sigma}^2$

$$\mathcal{X} : \mathbb{E}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x}; \mathbf{z})\right] = \boldsymbol{\mu}, \mathrm{Var}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x}; \mathbf{z})\right] = \boldsymbol{\sigma}^2. \tag{15}$$

Precisely, the emergent property statistics $f(\mathbf{x}; \mathbf{z})$ must have means $\boldsymbol{\mu}$ and variances $\boldsymbol{\sigma}^2$ over the EPI distribution of parameters and stochasticity of the data given the parameters.

In EPI, deep probability distributions are used as posterior approximations $q_{\boldsymbol{\theta}}(\mathbf{z} \mid \mathcal{X})$. In deep probability distributions, a simple random variable $\mathbf{z}_0 \sim q_0(\mathbf{z}_0)$ is mapped deterministically via a sequence of deep neural network layers $(g_1, \, .. \, g_l)$ parameterized by weights and biases $\boldsymbol{\theta}$ to the support of the distribution of interest:

$$\mathbf{z} = g_{\boldsymbol{\theta}}(\mathbf{z}_0) = g_l(..g_1(\mathbf{z}_0)) \sim q_{\boldsymbol{\theta}}(\mathbf{z}). \tag{16}$$

Such deep probability distributions embed the posterior distribution in a deep network. Once optimized, this deep network representation has remarkably useful properties: immediate posterior sampling, and immediate probability, gradient, and Hessian evaluation at any parameter choice.

Given a choice of model $p(\mathbf{x} \mid \mathbf{z})$ and emergent property of interest $\mathcal{X}$, $q_{\boldsymbol{\theta}}(\mathbf{z})$ is optimized via the neural network parameters $\boldsymbol{\theta}$ to find a maximally entropic distribution $q^*_{\boldsymbol{\theta}}$ within the deep variational family $\mathcal{Q}$ producing the emergent property $\mathcal{X}$:

$$\begin{aligned} q_{\boldsymbol{\theta}}(\mathbf{z} \mid \mathcal{X}) = q^*_{\boldsymbol{\theta}}(\mathbf{z}) = \underset{q_{\boldsymbol{\theta}} \in Q}{\operatorname{argmax}} \, H(q_{\boldsymbol{\theta}}(\mathbf{z})) \\ \text{s.t. } \mathcal{X} \quad : \quad \mathbb{E}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x};\mathbf{z})\right] = \boldsymbol{\mu}, \operatorname{Var}_{\mathbf{z},\mathbf{x}}\left[f(\mathbf{x};\mathbf{z})\right] = \boldsymbol{\sigma}^2. \end{aligned} \tag{17}$$

Entropy is chosen as the normative selection principle, since we want the posterior to only contain structure predicated by the emergent property [50, 51]. This choice of selection principle is also that of standard Bayesian inference, and we derive an exact relation between EPI and variational inference (see Section 5.1.5). However, a key difference is that variational inference and other Bayesian methods do not constrain the predictions of their inferred posteriors. This optimization is executed using the algorithm of Maximum Entropy Flow Networks (MEFNs) [44].

In the remainder of Section 5.1, we will explain the finer details and motivation of the EPI method. First, we explain related approaches and what EPI introduces to this domain (Section 5.1.1). Second, we describe the special class of deep probability distributions used in EPI called normalizing flows (Section 5.1.2). Next, we explain the constrained optimization technique used to solve Equation 17 (Section 5.1.3). Then, we demonstrate the details of this optimization in a toy example (Section 5.1.4). Finally, we establish the known relationship between maximum entropy distributions and exponential families (Section 5.1.5), which is used to explain the relation between EPI and variational inference (Section 5.1.6).

5.1.1 Related approaches

When Bayesian inference problems lack conjugacy, scientists use approximate inference methods like variational inference (VI) [73] and Markov chain Monte Carlo (MCMC) [74, 75]. After optimization, variational methods return a parameterized posterior distribution, which we can analyze. Also, the variational approximating distribution class is often chosen such that it permits fast sampling. In contrast MCMC methods only produce samples from the approximated posterior distribution. No parameterized distribution is estimated, and additional samples are always generated with the same sampling complexity. Inference in models defined by systems of differential has been

demonstrated with MCMC [76], although this approach requires tractable likelihoods. Advancements have leveraged structure in stochastic differential equation models to improve likelihood approximations, thus expanding the domain of applicable models [77].

Likelihood-free (or "simulation-based") inference (LFI) [78] is model parameter inference in the absence of a tractable likelihood function. The most prevalent approach to LFI is approximate Bayesian computation [79], in which satisfactory parameter samples are kept from random prior sampling according to a rejection heuristic. The obtained set of parameters do not have a probabilities, and further insight about the model must be gained from examination of the parameter set and their generated activity. Methodological advances to ABC methods have come through the use of Markov chain Monte Carlo (MCMC-ABC) [80] and sequential Monte Carlo (SMC-ABC) [27] sampling techniques. SMC-ABC is considered state-of-the-art ABC, yet this approach still struggles to scale in dimensionality (cf. Fig. 4). Furthermore, once a parameter set has been obtained by SMC-ABC from a finite set of particles, the SMC-ABC algorithm must be run again with a new population of initialized particles to obtain additional samples.

For scientific model analysis, we seek a posterior distribution exhibiting the properties of a well-chosen variational approximation: a parametric form conferring analytic calculations, and trivial sampling time. For this reason, ABC and MCMC techniques are unattractive, since they only produce a set of parameter samples and have unchanging sampling rate. EPI executes likelihood-free inference using the MEFN [44] algorithm using a deep variational posterior approximation. The deep neural network of EPI defines the parametric form of the posterior approximation. Furthermore, the EPI distribution is constrained to produce an emergent property. In other words, the summary statistics of the posterior predictive distribution are fixed to have certain first and second moments. EPI optimization is enabled using stochastic gradient techniques in the spirit of likelihood-free variational inference [41]. The analytic relationship between EPI and variational inference is explained in Secton 5.1.6.

We note that, during our preparation and early presentation of this work [81, 82], another work has arisen with broadly similar goals: bringing statistical inference to mechanistic models of neural circuits ([83, 84, 85]).We are encouraged by this general problem being recognized by others in the community, and we emphasize that these works offer complementary neuroscientific contributions (different theoretical models of focus) and use different technical methodologies (ours is built on our prior work [44], theirs similarly [86]).

The method EPI differs from SNPE in some key ways. SNPE belongs to a "sequential" class of

recently developed LFI methods in which two neural networks are used for posterior inference. This first neural network is a normalizing flow used to estimate the posterior $p(\mathbf{z} \mid \mathbf{x})$ (SNPE) or the likelihood $p(\mathbf{x} \mid \mathbf{z})$ (sequential neural likelihood (SNL [87])). A recent advance uses an unconstrained neural network to estimate the likelihood ratio (sequential neural ratio estimation (SNRE [88]). In SNL and SNRE, MCMC sampling techniques are used to obtain samples from the approximated posterior. This contrasts with EPI and SNPE, which afford a normalizing flow approximation to the posterior, which facilitates immediate measurements of sample probability, gradient, or Hessian for system analysis. The second neural network in this sequential class of methods is the amortizer. This network maps data $\mathbf{x}$ (or statistics $f(\mathbf{x}; \mathbf{z})$ or model parameters $\mathbf{z}$ to the weights and biases of the first neural network. These methods are optimized on a conditional density (or ratio) estimation objective on a sequentially adapting finite sample-based approximation to the posterior.

The approximating fidelity of the first neural network in sequential approaches is optimized to generalize across the entire distribution it is conditioned upon. This optimization towards generalization of sequential methods can reduce the accuracy at the singular posterior of interest. Whereas in EPI, the entire expressivity of the normalizing flow is dedicated to learning a single distribution as well as possible. While amortization is not possible in EPI parameterized by the mean parameter μ (due to the inverse mapping problem [89]), we have shown this two-network amortization approach to be effective in exponential family distributions defined by their natural parameterization [90].

Structural identifiability analysis involves the measurement of sensitivity and unidentifiabilities in natural models. Around a point, one can measure the Jacobian. One approach that scales well is EAR [91]. A popular efficient approach for systems of ODEs has been neural ODE adjoint [92] and its stochastic adaptation [93]. Casting identifiability as a statistical estimation problem, the profile likelihood can assess via iterated optimization while holding parameters fixed [94]. An exciting recent method is capable of recovering the functional form of such unidentifiabilities away from a point by following degenerate dimensions of the fisher information matrix [95]. Global structural non-identifiabilities can be found for models with polynomial or rational dynamics equations using DAISY [96]. With EPI, we have all the benefits given by a statistical inference method plus the ability to query the gradient or Hessian of the inferred distribution at any chosen parameter value.

5.1.2 Normalizing flows

Deep probability distributions are comprised of multiple layers of fully connected neural networks (Equation). When each neural network layer is restricted to be a bijective function, the sample density can be calculated using the change of variables formula at each layer of the network. For $\mathbf{z}_i = g_i(\mathbf{z_{i-1}})$,

$$p(\mathbf{z}_i) = p(g_i^{-1}(\mathbf{z}_i)) \left| \det \frac{\partial g_i^{-1}(\mathbf{z}_i)}{\partial \mathbf{z}_i} \right| = p(\mathbf{z}_{i-1}) \left| \det \frac{\partial g_i(\mathbf{z}_{i-1})}{\partial \mathbf{z}_{i-1}} \right|^{-1}. \tag{18}$$

However, this computation has cubic complexity in dimensionality for fully connected layers. By restricting our layers to normalizing flows [42, 97] – bijective functions with fast log determinant Jacobian computations, which confer a fast calculation of the sample log probability. Fast log probability calculation confers efficient optimization of the maximum entropy objective (see Section 5.1.3). We use the Real NVP [98] normalizing flow class, because its coupling architecture confers both fast sampling (forward) and fast log probability evaluation (backward). Fast probability evaluation in turn facilitates fast gradient and Hessian evaluation of log probability throughout parameter space. Glow permutations were used in between coupling stages [99]. This is in contrast to autoregressive architectures [43, 100], in which only forward or backward passes are efficient. Normalizing flow architectures for deep probability distributions used in EPI are specified by the number of coupling stages, neural network layers and units per layer for conditioning functions.

5.1.3 Augmented Lagrangian optimization

To optimize $q_{\boldsymbol{\theta}}(\mathbf{z})$ in Equation 17, the constrained optimization is executed using the augmented Lagrangian method. The following objective is minimized:

$$L(\boldsymbol{\theta}; \boldsymbol{\eta}_{\text{opt}}, c) = -H(q_{\boldsymbol{\theta}}) + \boldsymbol{\eta}_{\text{opt}}^{\top} R(\boldsymbol{\theta}) + \frac{c}{2} ||R(\boldsymbol{\theta})||^2 \tag{19}$$

where $R(\boldsymbol{\theta}) = \mathbb{E}_{\mathbf{z} \sim q_{\boldsymbol{\theta}}} \left[\mathbb{E}_{\mathbf{x} \sim p(\mathbf{x}|\mathbf{z})} \left[T(\mathbf{x}; \mathbf{z}) - \boldsymbol{\mu}_{\text{opt}} \right] \right]$, $\boldsymbol{\eta}_{\text{opt}} \in \mathbb{R}^m$ are the Lagrange multipliers where $m = |\boldsymbol{\mu}_{\text{opt}}| = |T(\mathbf{x}; \mathbf{z})|$, and c is the penalty coefficient. These Lagrange multipliers are closely related to the natural parameters $\boldsymbol{\eta}$ of exponential families (see Section 5.1.6). Deep neural network weights and biases $\boldsymbol{\theta}$ of the deep probability distribution are optimized according to Equation 19 using the Adam optimizer with learning rate 10^{-3} [101].

To take gradients with respect to the entropy $H(q_{\boldsymbol{\theta}}(\mathbf{z}))$, it can be expressed using the reparameterization trick as an expectation of the negative log density of parameter samples $\mathbf{z}$ over the

randomness in the parameterless initial distribution $q_0(\mathbf{z}_0)$:

$$H(q_{\boldsymbol{\theta}}(\mathbf{z})) = \int -q_{\boldsymbol{\theta}}(\mathbf{z}) \log(q_{\boldsymbol{\theta}}(\mathbf{z}))d\mathbf{z} = \mathbb{E}_{\mathbf{z}\sim q_{\boldsymbol{\theta}}} \left[-\log(q_{\boldsymbol{\theta}}(\mathbf{z}))\right] = \mathbb{E}_{\mathbf{z}_0\sim q_0} \left[-\log(q_{\boldsymbol{\theta}}(g_{\boldsymbol{\theta}}(\mathbf{z}_0)))\right]. \quad (20)$$

Thus, the gradient of the entropy of the deep probability distribution can be estimated as an average with respect to the base distribution $\mathbf{z}_0$:

$$\nabla_{\boldsymbol{\theta}} H(q_{\boldsymbol{\theta}}(\mathbf{z})) = \mathbb{E}_{\mathbf{z}_0\sim q_0} \left[-\nabla_{\boldsymbol{\theta}} \log(q_{\boldsymbol{\theta}}(g_{\boldsymbol{\theta}}(\mathbf{z}_0)))\right]. \quad (21)$$

The lagrangian parameters $\boldsymbol{\eta}_{\text{opt}}$ are initialized to zero and adapted following each augmented Lagrangian epoch, which is a period of optimization with fixed ($\boldsymbol{\eta}_{\text{opt}}$, c) for a given number of stochastic optimization iterations. A low value of c is used initially, and conditionally increased after each epoch based on constraint error reduction. The penalty coefficient is updated based on the result of a hypothesis test regarding the reduction in constraint violation. The p-value of $\mathbb{E}[||R(\boldsymbol{\theta}_{k+1})||] > \gamma \mathbb{E}\left[||R(\boldsymbol{\theta}_k)||\right]$ is computed, and c_{k+1} is updated to βc_k with probability $1-p$. The other update rule is $\boldsymbol{\eta}_{\text{opt},k+1} = \boldsymbol{\eta}_{\text{opt},k} + c_k \frac{1}{n} \sum_{i=1}^{n} (T(\mathbf{x}^{(i)}) - \boldsymbol{\mu}_{\text{opt}})$ given a batch size n. Throughout the study, $\beta = 4.0$, $\gamma = 0.25$, and the batch size was a hyperparameter, which varied according to the application of EPI.

The intention is that c and $\boldsymbol{\eta}_{\text{opt}}$ start at values encouraging entropic growth early in optimization. With each training epoch in which the update rule for c is invoked by unsatisfactory constraint error reduction, the constraint satisfaction terms are increasingly weighted, resulting in a decreased entropy. This encourages the discovery of suitable regions of parameter space, and the subsequent refinement of the distribution to produce the emergent property (see example in Section 5.1.4). The momentum parameters of the Adam optimizer are reset at the end of each augmented Lagrangian epoch.

Rather than starting optimization from some $\boldsymbol{\theta}$ drawn from a randomized distribution, we found that initializing $q_{\boldsymbol{\theta}}(\mathbf{z})$ to approximate an isotropic Gaussian distribution conferred more stable, consistent optimization. The parameters of the Gaussian initialization were chosen on an application-specific basis. Throughout the study, we chose isotropic Gaussian initializations with mean $\boldsymbol{\mu}_{\text{init}}$ at the center of the distribution support and some standard deviation $\boldsymbol{\sigma}_{\text{init}}$, except for one case, where an initialization informed by random search was used (see Section 5.2.1).

To assess whether the EPI distribution $q_{\boldsymbol{\theta}}(\mathbf{z})$ produces the emergent property, we defined a hypothesis testing convergence criteria. The algorithm has converged when a null hypothesis test of constraint violations $R(\boldsymbol{\theta})_i$ being zero is accepted for all constraints $i \in \{1, ..., m\}$ at a significance

threshold $\alpha = 0.05$. This significance threshold is adjusted through Bonferroni correction according to the number of constraints m. The p-values for each constraint are calculated according to a two-tailed nonparametric test, where 200 estimations of the sample mean $R(\boldsymbol{\theta})^i$ are made from k resamplings of $\mathbf{z}$ from a finite sample of size n taken at the end of the augmented Lagrangian epoch.

When assessing the suitability of EPI for a particular modeling question, there are some important technical considerations. First and foremost, as in any optimization problem, the defined emergent property should always be appropriately conditioned (constraints should not have wildly different units). Furthermore, if the program is underconstrained (not enough constraints), the distribution grows (in entropy) unstably unless mapped to a finite support. If overconstrained, there is no parameter set producing the emergent property, and EPI optimization will fail (appropriately). Next, one should consider the computational cost of the gradient calculations. In the best circumstance, there is a simple, closed form expression (e.g. Section 5.2.4) for the emergent property statistic given the model parameters. On the other end of the spectrum, many forward simulation iterations may be required before a high quality measurement of the emergent property statistic is available (e.g. Section 5.2.1). In such cases, backpropagating gradients through the SDE evolution will be expensive.

5.1.4 Example: 2D LDS

To gain intuition for EPI optimization, consider a two-dimensional linear dynamical system (2D LDS) model (Fig. 5A):

$$\tau \frac{d\mathbf{x}}{dt} = A\mathbf{x} \tag{22}$$

where

$$A = \begin{bmatrix} a_{1,1} & a_{1,2} \\ a_{2,1} & a_{2,2} \end{bmatrix}, \tag{23}$$

$\tau = 1$s, and $\mathbf{z} = [a_{1,1}, a_{1,2}, a_{2,1}, a_{2,2}]^\top$.

To learn the distribution of real entries of A that produce a band of oscillating systems around 1Hz, we formalized this emergent property as real(λ_1) having mean zero and the oscillation frequency 2πimag(λ_1) having mean $\omega = 1$ Hz. Thus, our statistics vector $f(\mathbf{x}; \mathbf{z})$ for this emergent property

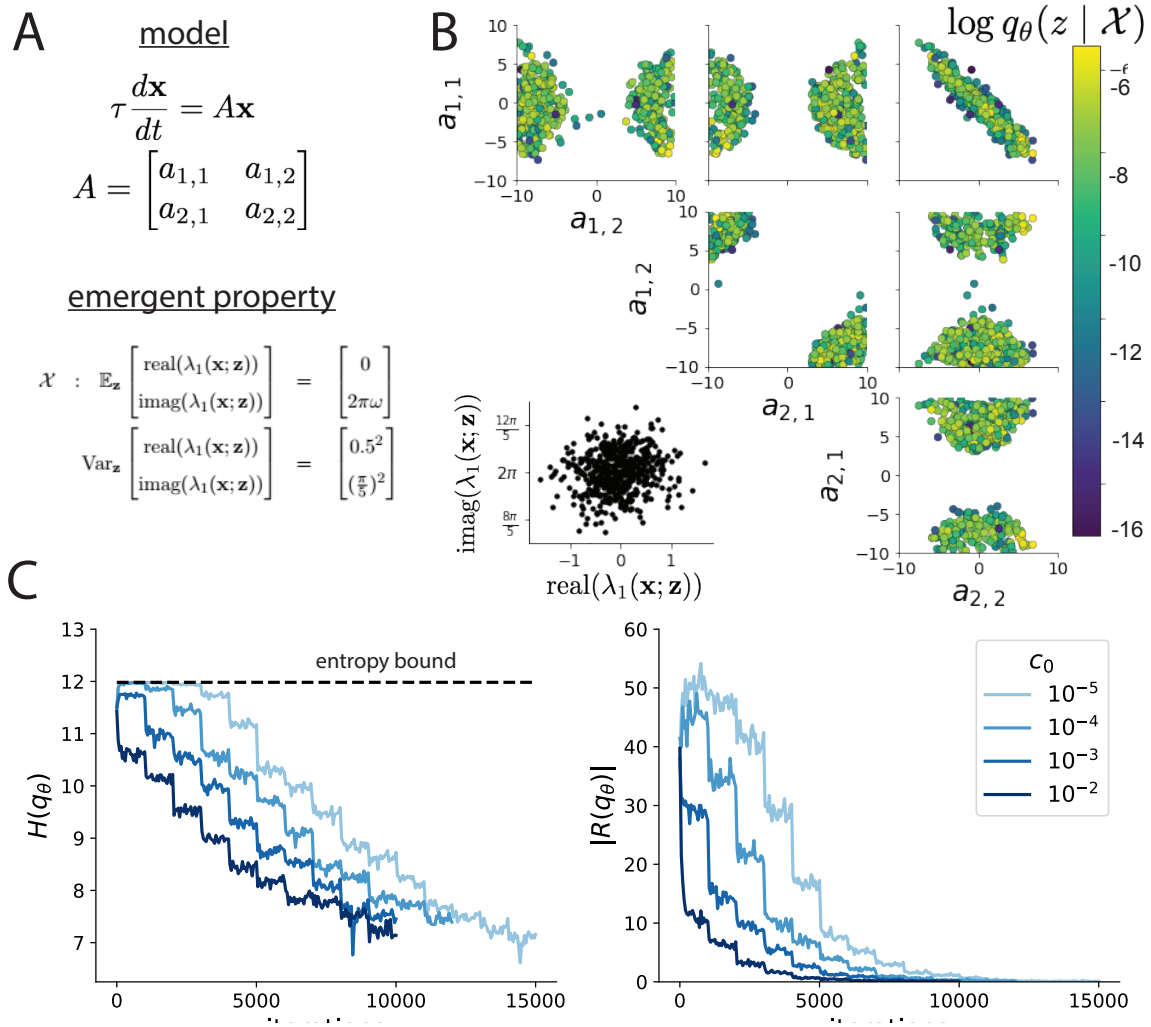

Figure 5: (LDS1): A. Two-dimensional linear dynamical system model, where real entries of the dynamics matrix A are the parameters. B. The EPI distribution for a two-dimensional linear dynamical system with $\tau = 1$ that produces oscillations around 1Hz. Inset show distribution of eigenvalues across the EPI posterior. C. (Left) Entropy throughout the optimization with different c_0. Individual optimizations were run with the same architecture and random seed. Norm of emergent property statistic errors throughout optimization.

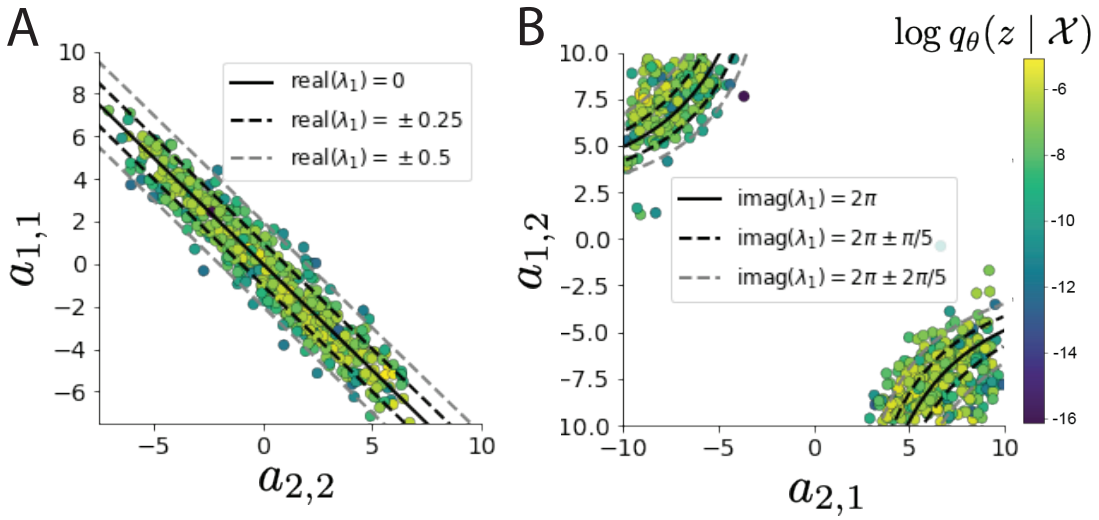

Figure 6: (LDS2): A. Probability contours in the $a_{1,1}$-$a_{2,2}$ plane were derived from the relationship to emergent property statistic of growth/decay factor real(λ_1). B. Probability contours in the $a_{1,2}$-$a_{2,1}$ plane were derived from the emergent property statistic of oscillation frequency 2πimag(λ_1).

is two dimensional, and after specifying the variance constraint is

$$\begin{aligned} \mathcal{X} \quad : \quad \mathbb{E}_{\mathbf{z}} \begin{bmatrix} \text{real}(\lambda_1(\mathbf{x};\mathbf{z})) \\ \text{imag}(\lambda_1(\mathbf{x};\mathbf{z})) \end{bmatrix} &= \begin{bmatrix} 0 \\ 2\pi\omega \end{bmatrix} \\ \text{Var}_{\mathbf{z}} \begin{bmatrix} \text{real}(\lambda_1(\mathbf{x};\mathbf{z})) \\ \text{imag}(\lambda_1(\mathbf{x};\mathbf{z})) \end{bmatrix} &= \begin{bmatrix} 0.5^2 \\ (\frac{\pi}{5})^2 \end{bmatrix}. \end{aligned} \tag{24}$$

Notation of mean and variance taken with respect to $\mathbf{x}$ is omitted, since this model is deterministic, and EPI generalizes naturally to this scenario.

Unlike the models we presented in the main text, this model admits an analytical form for the mean emergent property statistics given parameter $\mathbf{z}$, since the eigenvalues can be calculated using the quadratic formula:

$$\lambda = \frac{(\frac{a_{1,1}+a_{2,2}}{\tau}) \pm \sqrt{(\frac{a_{1,1}+a_{2,2}}{\tau})^2 + 4(\frac{a_{1,2}a_{2,1}-a_{1,1}a_{2,2}}{\tau})}}{2}. \tag{25}$$

Importantly, even though $f(\mathbf{x};\mathbf{z})$ is analytically available and does not require simulation, we cannot derive the distribution $q^*_{\boldsymbol{\theta}}(z \mid \mathcal{X})$ directly. This fact is due to the formally hard problem of the backward mapping: finding the natural parameters η from the mean parameters $\boldsymbol{\mu}$ of an exponential family distribution [89]. Instead, we can use EPI to approximate this distribution (Fig. 5B). We used a real-NVP normalizing flow architecture with three masks, two neural network layers of 50 units per mask, mapped onto a support of $z_i \in [-10, 10]$. (see Section 5.1.2).

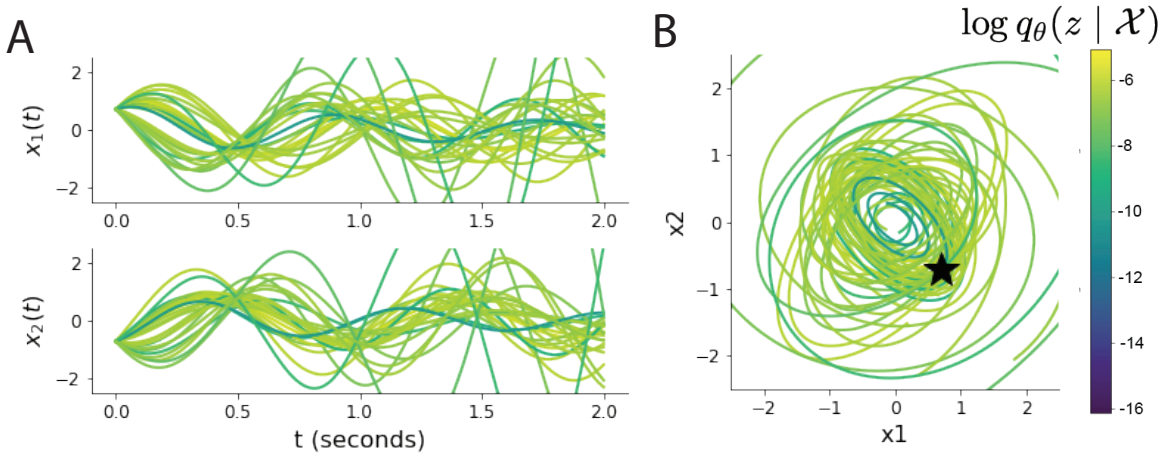

Figure 7: (LDS3): Sampled dynamical systems $\mathbf{z} \sim q_{\boldsymbol{\theta}}(\mathbf{z} \mid \mathcal{X})$ and their simulated activity from $\mathbf{x}(0) = [\frac{\sqrt{2}}{2}, -\frac{\sqrt{2}}{2}]$ colored by log probability. A. Each dimension of the simulated trajectories throughout time. B The simulated trajectories in phase space.

The effect of c_0 on the EPI optimization is shown in Figure 5C. Each EPI optimization was run for enough augmented Lagrangian epochs until convergence, with 1,000 iterations per epoch. We see that low values of c_0 (e.g. 10^{-5}) result in fast entropic growth to the uniform distribution at the outset of optimization. We know this, since the dashed black line indicates the maximum entropy of a distribution with the given bounds. Initial entropic growth can be desireable based on the application, however one should expect a greater number of epochs until convergence compared to a greater value of c_0 (e.g. 10^{-2}).

However, an important drawback of selecting a high 'c0' is that you may miss interesting structure in the learned distribution.

Even this relatively simple system has nontrivial (though intuitively sensible) structure in the parameter distribution. To validate our method, we analytically derived the contours of the probability density from the emergent property statistics and values. In the a_1-a_4 plane, the black line at $\text{real}(\lambda_1) = \frac{a_{1,1}+a_{2,2}}{2} = 0$, dotted black line at the standard deviation $\text{real}(\lambda_1) = \frac{a_{1,1}+a_{2,2}}{2} \pm 0.5$, and the dotted gray line at twice the standard deviation $\text{real}(\lambda_1) = \frac{a_{1,1}+a_{2,2}}{2} \pm 1$ follow the contour of probability density of the samples (Fig. 6A). The distribution precisely reflects the desired statistical constraints and model degeneracy in the sum of $a_{1,1}$ and $a_{2,2}$. Intuitively, the parameters equivalent with respect to emergent property statistic $\text{real}(\lambda_1)$ have similar log densities.

To explain the bimodality of the EPI distribution, we examined the imaginary component of λ_1.

When $\text{real}(\lambda_1) = \frac{a_{1,1}+a_{2,2}}{2} = 0$, we have

$$\text{imag}(\lambda_1) = \begin{cases} \sqrt{\frac{a_{1,1}a_{2,2}-a_{1,2}a_{2,1}}{\tau}}, & \text{if } a_{1,1}a_{2,2} < a_{1,2}a_{2,1} \\ 0 & \text{otherwise} \end{cases}. \tag{26}$$

When $\tau = 1$ and $a_{1,1}a_{2,2} > a_{1,2}a_{2,1}$ (center of distribution above), we have the following equation for the other two dimensions:

$$\text{imag}(\lambda_1)^2 = a_{1,1}a_{2,2} - a_{1,2}a_{2,1} \tag{27}$$

Since we constrained $\mathbb{E}_{\mathbf{z}\sim q_{\boldsymbol{\theta}}}[\text{imag}(\lambda)] = 2\pi$ (with $\omega = 1$), we can plot contours of the equation $\text{imag}(\lambda_1)^2 = a_{1,1}a_{2,2} - a_{1,2}a_{2,1} = (2\pi)^2$ assuming $a_{1,1}a_{2,2} = \mathbb{E}_z[a_{1,1}a_{2,2}]$ (Fig. 6B). This validates the curved structure of the inferred distribution learned through EPI. Subtler combinations of model and emergent property will have more complexity, further motivating the use of EPI for understanding these systems. As we expect, the distribution results in samples of two-dimensional linear systems oscillating near 1Hz (Fig. 7).

5.1.5 Maximum entropy distributions and exponential families

Maximum entropy distributions have a fundamental link to exponential family distributions. A maximum entropy distribution of form:

$$\begin{aligned} p^*(\mathbf{z}) &= \underset{p\in\mathcal{P}}{\text{argmax}}\, H(p(\mathbf{z})) \\ &\text{s.t. } \mathbb{E}_{\mathbf{z}\sim p}[T(\mathbf{z})] = \boldsymbol{\mu}_{\text{opt}}. \end{aligned} \tag{28}$$

will have probability density in the exponential family:

$$p^*(\mathbf{z}) \propto \exp(\boldsymbol{\eta}^\top T(\mathbf{z})). \tag{29}$$

The mappings between the mean parameterization $\boldsymbol{\mu}_{\text{opt}}$ and the natural parameterization $\boldsymbol{\eta}$ are formally hard to identify [89].

In EPI, emergent properties are defined as statistics having a fixed mean and variance as in Equation 2

$$\mathbb{E}_{\mathbf{z},\mathbf{x}}[f(\mathbf{x};\mathbf{z})] = \boldsymbol{\mu}, \text{Var}_{\mathbf{z},\mathbf{x}}[f(\mathbf{x};\mathbf{z})] = \boldsymbol{\sigma}^2. \tag{30}$$

The variance constraint is a second moment constraint on $f(\mathbf{x};\mathbf{z})$

$$\text{Var}_{\mathbf{z},\mathbf{x}}[f(\mathbf{x};\mathbf{z})] = \mathbb{E}_{\mathbf{z},\mathbf{x}}\left[(f(\mathbf{x};\mathbf{z}) - \boldsymbol{\mu})^2\right] \tag{31}$$

As a general maximum entropy distribution (Equation 28), the sufficient statistics vector contains both first and second order moments of $f(\mathbf{x}; \mathbf{z})$

$$T(\mathbf{x}; \mathbf{z}) = \begin{bmatrix} f(\mathbf{x}; \mathbf{z}) \\ (f(\mathbf{x}; \mathbf{z}) - \boldsymbol{\mu})^2 \end{bmatrix}, \tag{32}$$

which are constrained to the chosen means and variances

$$\boldsymbol{\mu}_{\text{opt}} = \begin{bmatrix} \boldsymbol{\mu} \\ \boldsymbol{\sigma}^2 \end{bmatrix}. \tag{33}$$

5.1.6 EPI as variational inference

In Bayesian inference a prior belief about model parameters $\mathbf{z}$ is stated in a prior distribution $p(\mathbf{z})$, and the statistical model capturing the effect of $\mathbf{z}$ on observed data points $\mathbf{x}$ is formalized in the likelihood distribution $p(\mathbf{x} \mid \mathbf{z})$. In Bayesian inference, we obtain a posterior distribution $p(z \mid \mathbf{x})$, which captures how the data inform our knowledge of model parameters using Bayes' rule:

$$p(\mathbf{z} \mid \mathbf{x}) = \frac{p(\mathbf{x} \mid \mathbf{z})p(\mathbf{z})}{p(\mathbf{x})}. \tag{34}$$

The posterior distribution is analytically available when the prior is conjugate with the likelihood. However, conjugacy is rare in practice, and alternative methods, such as variational inference [102], are utilized.

In variational inference, a posterior approximation $q^*_{\boldsymbol{\theta}}$ is chosen from within some variational family $\mathcal{Q}$

$$q^*_{\boldsymbol{\theta}}(\mathbf{z}) = \underset{q_{\boldsymbol{\theta}} \in Q}{\operatorname{argmin}} \, KL(q_{\boldsymbol{\theta}}(\mathbf{z}) \mid\mid p(\mathbf{z} \mid \mathbf{x})). \tag{35}$$

The KL divergence can be written in terms of entropy of the variational approximation:

$$KL(q_{\boldsymbol{\theta}}(\mathbf{z}) \mid\mid p(\mathbf{z} \mid \mathbf{x})) = \mathbb{E}_{\mathbf{z} \sim q_{\boldsymbol{\theta}}} \left[\log(q_{\boldsymbol{\theta}}(\mathbf{z}))\right] - \mathbb{E}_{\mathbf{z} \sim q_{\boldsymbol{\theta}}} \left[\log(p(\mathbf{z} \mid \mathbf{x}))\right] \tag{36}$$

$$= -H(q_{\boldsymbol{\theta}}) - \mathbb{E}_{\mathbf{z} \sim q_{\boldsymbol{\theta}}} \left[\log(p(\mathbf{x} \mid \mathbf{z})) + \log(p(\mathbf{z})) - \log(p(\mathbf{x}))\right] \tag{37}$$

Since the marginal distribution of the data $p(\mathbf{x})$ (or "evidence") is independent of $\boldsymbol{\theta}$, variational inference is executed by optimizing the remaining expression. This is usually framed as maximizing the evidence lower bound (ELBO)

$$\underset{q_{\boldsymbol{\theta}} \in Q}{\operatorname{argmin}} \, KL(q_{\boldsymbol{\theta}} \mid\mid p(\mathbf{z} \mid \mathbf{x})) = \underset{q_{\boldsymbol{\theta}} \in Q}{\operatorname{argmax}} \, H(q_{\boldsymbol{\theta}}) + \mathbb{E}_{\mathbf{z} \sim q_{\boldsymbol{\theta}}} \left[\log(p(\mathbf{x} \mid \mathbf{z})) + \log(p(\mathbf{z}))\right]. \tag{38}$$

Now, consider the setting where we have chosen a uniform prior, and stipulate a mean-field gaussian likelihood on a chosen statistic of the data $f(\mathbf{x}; \mathbf{z})$

$$p(\mathbf{x} \mid \mathbf{z}) = \mathcal{N}(f(\mathbf{x}; \mathbf{z}) \mid \boldsymbol{\mu}_f, \Sigma_f), \tag{39}$$

where $\Sigma_f = \text{diag}(\boldsymbol{\sigma}_f^2)$. The log likelihood is then proportional to a dot product of the natural parameter of this mean-field gaussian distribution and the first and second moment statistics.

$$\log p(\mathbf{x} \mid \mathbf{z}) \propto \boldsymbol{\eta}_f^\top T(\mathbf{x}, \mathbf{z}), \tag{40}$$

where

$$\boldsymbol{\eta}_f = \begin{bmatrix} \frac{\boldsymbol{\mu}_f}{\boldsymbol{\sigma}_f^2} \\ \frac{-1}{2\boldsymbol{\sigma}_f^2} \end{bmatrix}, \text{ and} \tag{41}$$

$$T(\mathbf{x}; \mathbf{z}) = \begin{bmatrix} f(\mathbf{x}; \mathbf{z}) \\ (f(\mathbf{x}; \mathbf{z}) - \boldsymbol{\mu}_f)^2 \end{bmatrix}. \tag{42}$$

The variational objective is then

$$\underset{q_{\boldsymbol{\theta}} \in Q}{\text{argmax}}\, H(q_{\boldsymbol{\theta}}) + \boldsymbol{\eta}_f^\top \mathbb{E}_{\mathbf{z} \sim q_{\boldsymbol{\theta}}} \left[T(\mathbf{x}; \mathbf{z}) \right] \tag{43}$$

Comparing this to the Lagrangian objective (without augmentation) of EPI, we see they are the same

$$\begin{aligned} q^*_{\boldsymbol{\theta}}(\mathbf{z}) &= \underset{q_{\boldsymbol{\theta}} \in Q}{\text{argmin}} -H(q_{\boldsymbol{\theta}}) + \boldsymbol{\eta}_{\text{opt}}^\top \left(\mathbb{E}_{\mathbf{z},\mathbf{x}} \left[T(\mathbf{x}; \mathbf{z}) \right] - \boldsymbol{\mu}_{\text{opt}} \right) \\ &= \underset{q_{\boldsymbol{\theta}} \in Q}{\text{argmin}} -H(q_{\boldsymbol{\theta}}) + \boldsymbol{\eta}_{\text{opt}}^\top \mathbb{E}_{\mathbf{z},\mathbf{x}} \left[T(\mathbf{x}; \mathbf{z}) \right]. \end{aligned} \tag{44}$$

where $T(\mathbf{x}; \mathbf{z})$ consists of the first and second moments of the emergent property statistic $f(\mathbf{x}; \mathbf{z})$ (Equation 32). Thus, EPI is implicitly executing variational inference with a uniform prior and a mean-field gaussian likelihood on the emergent property statistics. The data $\mathbf{x}$ used by this implicit variational inference program would be that generated by the adapting variational approximation $\mathbf{x} \sim p(\mathbf{x} \mid \mathbf{z}) q_{\boldsymbol{\theta}}(\mathbf{z})$, and the likelihood parameters $\boldsymbol{\eta}_f$ of EPI optimization epoch k are predicated by $\boldsymbol{\eta}_{\text{opt},k}$. However, in EPI we have not specified a prior distribution, or collected data, which can inform us about model parameters. Instead we have a mathematical specification of an emergent property, which the model must produce, and a maximum entropy selection principle. Accordingly, we replace the notation of $p(\mathbf{z} \mid \mathbf{x})$ with $p(\mathbf{z} \mid \mathcal{X})$ conceptualizing an inferred distribution that obeys emergent property $\mathcal{X}$ (see Section 5.1).

5.2 Theoretical models

In this study, we used emergent property inference to examine several models relevant to theoretical neuroscience. Here, we provide the details of each model and the related analyses.

5.2.1 Stomatogastric ganglion

We analyze how the parameters $\mathbf{z} = [g_{\text{el}}, g_{\text{synA}}]$ govern the emergent phenomena of intermediate hub frequency in a model of the stomatogastric ganglion (STG) [49] shown in Figure 1A with activity $\mathbf{x} = [x_{\text{f1}}, x_{\text{f2}}, x_{\text{hub}}, x_{\text{s1}}, x_{\text{s2}}]$, using the same hyperparameter choices as Gutierrez et al. Each neuron's membrane potential $x_\alpha(t)$ for $\alpha \in \{\text{f1}, \text{f2}, \text{hub}, \text{s1}, \text{s2}\}$ is the solution of the following stochastic differential equation:

$$C_m \frac{dx_\alpha}{dt} = -\left[h_{leak}(\mathbf{x};\mathbf{z}) + h_{Ca}(\mathbf{x};\mathbf{z}) + h_K(\mathbf{x};\mathbf{z}) + h_{hyp}(\mathbf{x};\mathbf{z}) + h_{elec}(\mathbf{x};\mathbf{z}) + h_{syn}(\mathbf{x};\mathbf{z})\right] + dB. \quad (45)$$

The input current of each neuron is the sum of the leak, calcium, potassium, hyperpolarization, electrical and synaptic currents as well as gaussian noise dB. Each current component is a function of all membrane potentials and the conductance parameters $\mathbf{z}$.

The capacitance of the cell membrane was set to $C_m = 1nF$. Specifically, the currents are the difference in the neuron's membrane potential and that current type's reversal potential multiplied by a conductance:

$$h_{leak}(\mathbf{x};\mathbf{z}) = g_{leak}(x_\alpha - V_{leak}) \quad (46)$$

$$h_{elec}(\mathbf{x};\mathbf{z}) = g_{\text{el}}(x_\alpha^{post} - x_\alpha^{pre}) \quad (47)$$

$$h_{syn}(\mathbf{x};\mathbf{z}) = g_{syn} S_\infty^{pre}(x_\alpha^{post} - V_{syn}) \quad (48)$$

$$h_{Ca}(\mathbf{x};\mathbf{z}) = g_{Ca} M_\infty(x_\alpha - V_{Ca}) \quad (49)$$

$$h_K(\mathbf{x};\mathbf{z}) = g_K N(x_\alpha - V_K) \quad (50)$$

$$h_{hyp}(\mathbf{x};\mathbf{z}) = g_h H(x_\alpha - V_{hyp}). \quad (51)$$

The reversal potentials were set to $V_{leak} = -40mV$, $V_{Ca} = 100mV$, $V_K = -80mV$, $V_{hyp} = -20mV$, and $V_{syn} = -75mV$. The other conductance parameters were fixed to $g_{leak} = 1 \times 10^{-4}\mu S$. g_{Ca}, g_K, and g_{hyp} had different values based on fast, intermediate (hub) or slow neuron. The fast conductances had values $g_{Ca} = 1.9\times10^{-2}$, $g_K = 3.9\times10^{-2}$, and $g_{hyp} = 2.5\times10^{-2}$. The intermediate conductances had values $g_{Ca} = 1.7 \times 10^{-2}$, $g_K = 1.9 \times 10^{-2}$, and $g_{hyp} = 8.0 \times 10^{-3}$. Finally, the slow conductances had values $g_{Ca} = 8.5 \times 10^{-3}$, $g_K = 1.5 \times 10^{-2}$, and $g_{hyp} = 1.0 \times 10^{-2}$.

Furthermore, the Calcium, Potassium, and hyperpolarization channels have time-dependent gating dynamics dependent on steady-state gating variables M_∞, N_∞ and H_∞, respectively:

$$M_\infty = 0.5\left(1 + \tanh\left(\frac{x_\alpha - v_1}{v_2}\right)\right) \tag{52}$$

$$\frac{dN}{dt} = \lambda_N(N_\infty - N) \tag{53}$$

$$N_\infty = 0.5\left(1 + \tanh\left(\frac{x_\alpha - v_3}{v_4}\right)\right) \tag{54}$$

$$\lambda_N = \phi_N \cosh\left(\frac{x_\alpha - v_3}{2v_4}\right) \tag{55}$$

$$\frac{dH}{dt} = \frac{(H_\infty - H)}{\tau_h} \tag{56}$$

$$H_\infty = \frac{1}{1 + \exp\left(\frac{x_\alpha + v_5}{v_6}\right)} \tag{57}$$

$$\tau_h = 272 - \left(\frac{-1499}{1 + \exp\left(\frac{-x_\alpha + v_7}{v_8}\right)}\right). \tag{58}$$

where we set $v_1 = 0mV$, $v_2 = 20mV$, $v_3 = 0mV$, $v_4 = 15mV$, $v_5 = 78.3mV$, $v_6 = 10.5mV$, $v_7 = -42.2mV$, $v_8 = 87.3mV$, $v_9 = 5mV$, and $v_{th} = -25mV$.

Finally, there is a synaptic gating variable as well:

$$S_\infty = \frac{1}{1 + \exp\left(\frac{v_{th} - x_\alpha}{v_9}\right)}. \tag{59}$$

When the dynamic gating variables are considered, this is actually a 15-dimensional nonlinear dynamical system. Gaussian noise of variance $(1 \times 10^{-12})^2$ amps makes the model stochastic, and introduces variability in frequency at each parameterization $\mathbf{z}$.

In order to measure the frequency of the hub neuron during EPI, the STG model was simulated for $T = 300$ time steps of $dt = 25ms$. The chosen dt and T were the most computationally convenient choices yielding accurate frequency measurement. We used a basis of complex exponentials with frequencies from 0.0-1.0 Hz at 0.01Hz resolution to measure frequency from simulated time series

$$\Phi = [0.0, 0.01, ..., 1.0]^\top .. \tag{60}$$

To measure spiking frequency, we processed simulated membrane potentials with a relu (spike extraction) and low-pass filter with averaging window of size 20, then took the frequency with the maximum absolute value of the complex exponential basis coefficients of the processed time-series. The first 20 temporal samples of the simulation are ignored to account for initial transients.

To differentiate through the maximum frequency identification, we used a soft-argmax Let $X_\alpha \in \mathcal{C}^{|\Phi|}$ be the complex exponential filter bank dot products with the signal $x_\alpha \in \mathbb{R}^N$, where $\alpha \in \{\text{f1}, \text{f2}, \text{hub}, \text{s1}, \text{s2}\}$. The soft-argmax is then calculated using temperature parameter $\beta = 100$

$$\psi_\alpha = \text{softmax}(\beta |X_\alpha| \odot i), \tag{61}$$

where $i = [0, 1, ..., 100]$. The frequency is then calculated as

$$\omega_\alpha = 0.01 \psi_\alpha \text{Hz}. \tag{62}$$

Intermediate hub frequency, like all other emergent properties in this work, is defined by the mean and variance of the emergent property statistics. In this case, we have one statistic, hub neuron frequency, where the mean was chosen to be 0.55Hz, and variance was chosen to be $(0.025\text{Hz})^2$ to capture variation in frequency between 0.5Hz and 0.6Hz (Equation 2). As a maximum entropy distribution, $T(\mathbf{x}, \mathbf{z})$ is comprised of both these first and second moments of the hub neuron frequency (as in Equations 32 and 33)

$$T(\mathbf{x}; \mathbf{z}) = \begin{bmatrix} \omega_{\text{hub}}(\mathbf{x}; \mathbf{z}) \\ (\omega_{\text{hub}}(\mathbf{x}; \mathbf{z}) - 0.55)^2 \end{bmatrix}, \tag{63}$$

$$\boldsymbol{\mu}_{\text{opt}} = \begin{bmatrix} 0.55 \\ 0.025^2 \end{bmatrix}. \tag{64}$$

Throughout optimization, the augmented Lagrangian parameters η and c, were updated after each epoch of 5,000 iterations(see Section 5.1.3). The optimization converged after five epochs (Fig. S4).

For EPI in Fig 1E, we used a real NVP architecture with three coupling layers of affine transformations parameterized by two-layer neural networks of 25 units per layer. The initial distribution was a standard isotropic gaussian $z_0 \sim \mathcal{N}(\mathbf{0}, I)$ mapped to a support of $\mathbf{z} = [g_{\text{el}}, g_{\text{synA}}] \in [4, 8] \times [0.01, 4]$. We did not include $g_{\text{synA}} < 0.01$, since conductances that low make the circuit simulations numerically unstable. We used an augmented Lagrangian coefficient of $c_0 = 10^5$, a batch size $n = 400$, set $\nu = 0.25$, and initialized $q_{\boldsymbol{\theta}}(\mathbf{z})$ to produce a gaussian approximation to samples returned from an initial ABC search. This initialization had much greater entropy and a different emergent property than the returned EPI posterior.

TODO write about specifics of the Hessian analysis.

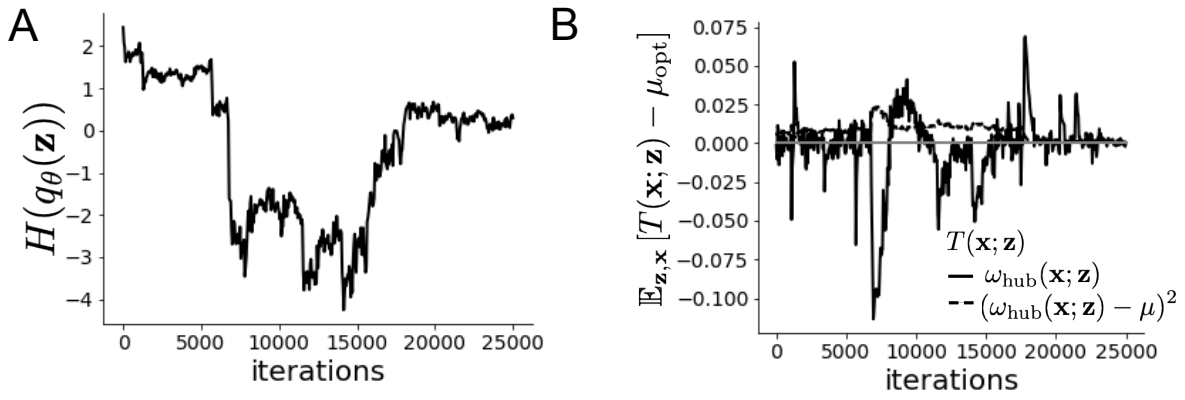

Figure 8: (STG1): EPI optimization of the STG model producing network syncing. A. Entropy throughout optimization. B. The emergent property statistic means and variances converge to their constraints at 25,000 iterations following the fifth augmented Lagrangian epoch.

5.2.2 Primary visual cortex

Connectivity (W_{fit}) and input ($\mathbf{h}_{b,\text{fit}}$ and $\mathbf{h}_{c,\text{fit}}$) parameters were fit using the deterministic V1 circuit model [64]

$$W_{\text{fit}} = \begin{bmatrix} W_{EE} & W_{EP} & W_{ES} & W_{EV} \\ W_{PE} & W_{PP} & W_{PS} & W_{PV} \\ W_{SE} & W_{SP} & W_{SS} & W_{SV} \\ W_{VE} & W_{VP} & W_{VS} & W_{VV} \end{bmatrix} = \begin{bmatrix} 2.18 & -1.19 & -.594 & -.229 \\ 1.66 & -.651 & -.680 & -.242 \\ .895 & -5.22 \times 10^{-3} & -1.51 \times 10^{-4} & -.761 \\ 3.34 & -2.31 & -.254 & -2.52 \times 10^{-4} \end{bmatrix}, \tag{65}$$

$$\mathbf{h}_{b,\text{fit}} = \begin{bmatrix} .416 \\ .429 \\ .491 \\ .486 \end{bmatrix}, \tag{66}$$

and

$$\mathbf{h}_{c,\text{fit}} = \begin{bmatrix} .359 \\ .403 \\ 0 \\ 0 \end{bmatrix}. \tag{67}$$

To obtain rates on a realistic scale (100-fold greater), we map these fitted parameters to an equiv-

alence class

$$W = \begin{bmatrix} W_{EE} & W_{EP} & W_{ES} & W_{EV} \\ W_{PE} & W_{PP} & W_{PS} & W_{PV} \\ W_{SE} & W_{SP} & W_{SS} & W_{SV} \\ W_{VE} & W_{VP} & W_{VS} & W_{VV} \end{bmatrix} = \begin{bmatrix} .218 & -.119 & -.0594 & -.0229 \\ .166 & -.0651 & -.068 & -.0242 \\ .0895 & -5.22 \times 10^{-4} & -1.51 \times 10^{-5} & -.0761 \\ .334 & -.231 & -.0254 & -2.52 \times 10^{-5} \end{bmatrix}, \tag{68}$$

$$\mathbf{h}_b = \begin{bmatrix} h_{b,E} \\ h_{b,P} \\ h_{b,S} \\ h_{b,V} \end{bmatrix} = \begin{bmatrix} 4.16 \\ 4.29 \\ 4.91 \\ 4.86 \end{bmatrix}, \tag{69}$$

and

$$\mathbf{h}_c = \begin{bmatrix} h_{c,E} \\ h_{c,P} \\ h_{c,S} \\ h_{c,V} \end{bmatrix} = \begin{bmatrix} 3.59 \\ 4.03 \\ 0 \\ 0 \end{bmatrix}. \tag{70}$$

Since the E-population of this network increases exponentially in the absense of recurrent inhibitory feedback, we may also observe a paradoxical effect in the inhibitory populations (which is present in E-I networks). At 50% contrast (Fig. 2B, dots), this network exhibits a paradoxical effect in the P-population (Fig. 2C), but no others (Fig. 9). That is, for a small increase in h_P, $\mathbb{E}_t[x_P]$ decreases.

Fano factor is calculated as the temporal variance divided by the temporal mean following sometime t_{ss} following dynamical evolution from the initial state at $\mathbf{x}(t=0)$.

5.2.3 Superior colliculus

In the model of Duan et al [47], there are four total units: two in each hemisphere corresponding to the Pro/Contra and Anti/Ipsi populations. They are denoted as left Pro (LP), left Anti (LA), right Pro (RP) and right Anti (RA). Each unit has an activity (x_α) and internal variable (u_α) related by

$$x_\alpha = \phi(u_\alpha) = \left(\frac{1}{2}\tanh\left(\frac{u_\alpha - a}{b}\right) + \frac{1}{2}\right) \tag{71}$$

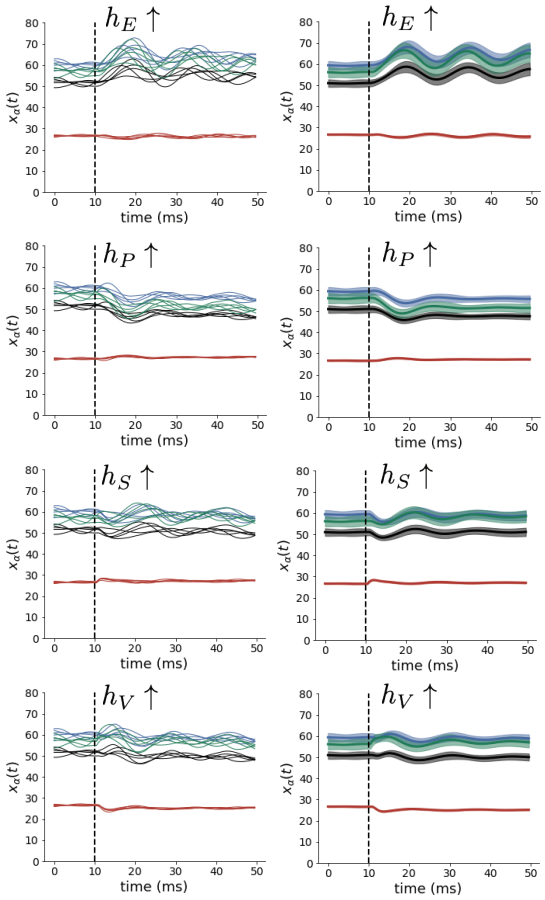

Figure 9: Supplemental Figure: (Left) Simulations for small increases in neuron-type population input. Input magnitudes are chosen so that effect is salient (0.002 for E and P, but 0.02 for S and V). (Right) Average and standard deviation of stochastic fluctuations of responses.

where $\alpha \in \{LP, LA, RA, RP\}$, $a = 0.05$ and $b = 0.5$ control the position and shape of the nonlinearity, respectively. During periods of optogenetic inactivation, activity was decreased proportional to the optogenetic strength γ

$$x_\alpha = (1 - \gamma)\phi(u_\alpha). \tag{72}$$

We order the neural populations of x and u in the following manner

$$\mathbf{x} = \begin{bmatrix} x_{LP} \\ x_{LA} \\ x_{RP} \\ x_{RA} \end{bmatrix} \qquad \mathbf{u} = \begin{bmatrix} u_{LP} \\ u_{LA} \\ u_{RP} \\ u_{RA} \end{bmatrix}, \tag{73}$$

which evolve according to

$$\tau \frac{d\mathbf{u}}{dt} = -\mathbf{u} + W\mathbf{x} + \mathbf{h} + d\mathbf{B}. \tag{74}$$

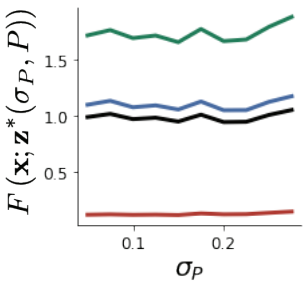

Figure 10: Supplemental Figure: Fano factors along the ridge of the posterior in Fig. 2E.

with time constant $\tau = 0.09s$, step size 24ms and Gaussian noise $d\mathbf{B}$ of variance 0.2. The weight matrix has 4 parameters sW, vW, hW, and dW:

$$W = \begin{bmatrix} sW & vW & hW & dW \\ vW & sW & dW & hW \\ hW & dW & sW & vW \\ dW & hW & vW & sW \end{bmatrix}. \tag{75}$$

The circuit receives four different inputs throughout each trial, which has a total length of 1.8s.

$$\mathbf{h} = \mathbf{h}_{\text{constant}} + \mathbf{h}_{\text{P,bias}} + \mathbf{h}_{\text{rule}} + \mathbf{h}_{\text{choice-period}} + \mathbf{h}_{\text{light}}. \tag{76}$$

There is a constant input to every population,

$$\mathbf{h}_{\text{constant}} = I_{\text{constant}}[1, 1, 1, 1]\top, \tag{77}$$

a bias to the Pro populations

$$\mathbf{h}_{\text{P,bias}} = I_{\text{P,bias}}[1, 0, 1, 0]\top, \tag{78}$$

rule-based input depending on the condition

$$\mathbf{h}_{\text{P,rule}}(t) = \begin{cases} I_{\text{P,rule}}[1, 0, 1, 0]^\top, & \text{if } t \leq 1.2s \\ 0, & \text{otherwise} \end{cases} \tag{79}$$

$$\mathbf{h}_{\text{A,rule}}(t) = \begin{cases} I_{\text{A,rule}}[0, 1, 0, 1]^\top, & \text{if } t \leq 1.2s \\ 0, & \text{otherwise} \end{cases}, \tag{80}$$

a choice-period input

$$\mathbf{h}_{\text{choice}}(t) = \begin{cases} I_{\text{choice}}[1, 1, 1, 1]^\top, & \text{if } t > 1.2s \\ 0, & \text{otherwise} \end{cases}, \tag{81}$$

and an input to the right or left-side depending on where the light stimulus is delivered

$$\mathbf{h}_{\text{light}}(t) = \begin{cases} I_{\text{light}}[1,1,0,0]^\top, & \text{if } 1.2s < t < 1.5s \text{ and Left} \\ I_{\text{light}}[0,0,1,1]^\top, & \text{if } 1.2s < t < 1.5s \text{ and Right} \\ 0, & \text{otherwise} \end{cases}. \tag{82}$$

The input parameterization was fixed to $I_{\text{constant}} = 0.75$, $I_{\text{P,bias}} = 0.5$, $I_{\text{P,rule}} = 0.6$, $I_{\text{A,rule}} = 0.6$, $I_{\text{choice}} = 0.25$, and $I_{\text{light}} = 0.5$.

The accuracies of p_P and p_A are calculated as

$$p_P(\mathbf{x};\mathbf{z}) = \mathbb{E}_{\mathbf{x}\sim p(\mathbf{x}|\mathbf{z})}\left[\Theta[x_{LP}(t = 1.8s) - x_{RP}(t = 1.8s)]\right] \tag{83}$$

and

$$p_A(\mathbf{x};\mathbf{z}) = \mathbb{E}_{\mathbf{x}\sim p(\mathbf{x}|\mathbf{z})}\left[\Theta[x_{RP}(t = 1.8s) - x_{LP}(t = 1.8s)]\right] \tag{84}$$

given that the stimulus is on the left side, where Θ is the Heaviside step function.

The Heaviside step function is approximated as

$$\Theta(\mathbf{x}) = \text{sigmoid}(\beta\mathbf{x}), \tag{85}$$

where $\beta = 100$.

As a maximum entropy distribution, $T(\mathbf{x},\mathbf{z})$ is comprised of both these first and second moments of the accuracy in each task (as in Equations 32 and 33)

$$T(\mathbf{x};\mathbf{z}) = \begin{bmatrix} p(\mathbf{x};\mathbf{z})_P \\ p(\mathbf{x};\mathbf{z})_A \\ (p(\mathbf{x};\mathbf{z})_P - 75\%)^2 \\ (p(\mathbf{x};\mathbf{z})_A - 75\%)^2 \end{bmatrix}, \tag{86}$$

$$\boldsymbol{\mu}_{\text{opt}} = \begin{bmatrix} 75\% \\ 75\% \\ 5\%^2 \\ 5\%^2 \end{bmatrix}. \tag{87}$$

Throughout optimization, the augmented Lagrangian parameters η and c, were updated after each epoch of 2,000 iterations(see Section 5.1.3). The optimization converged after six epochs (Fig. 15).

For EPI in Fig. 3C, we used a real NVP architecture with three coupling layers of affine transformations parameterized by two-layer neural networks of 50 units per layer. The initial distribution was

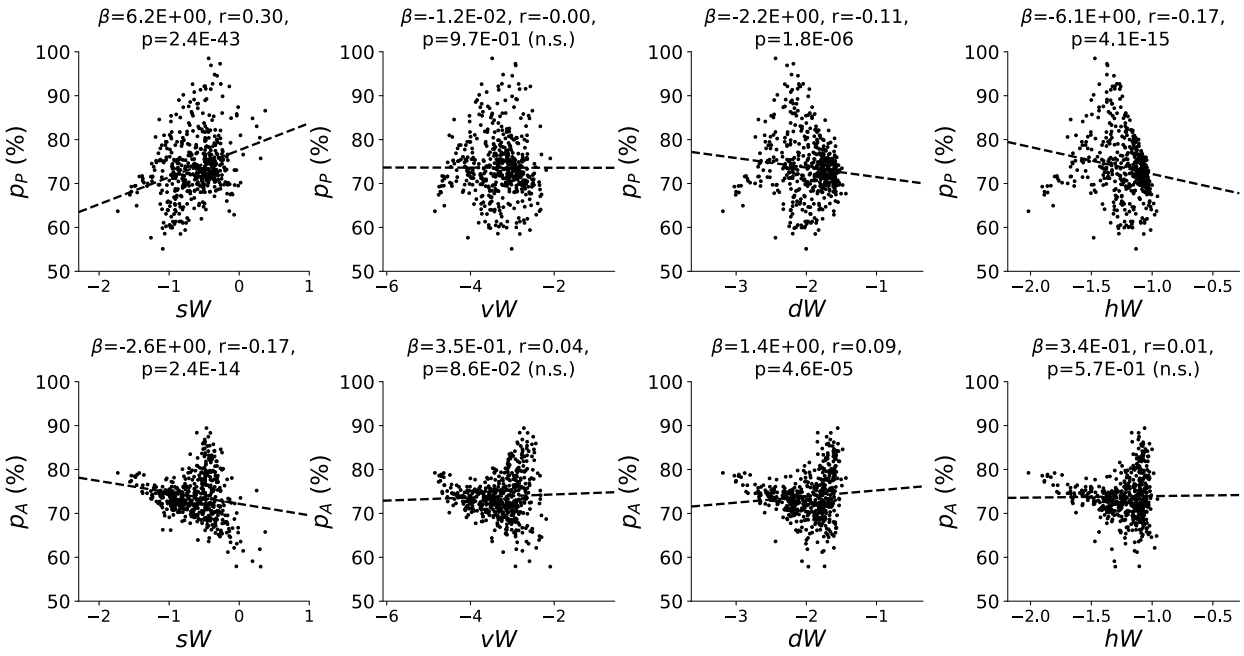

Figure 11: (SC1): Connectivity parameters of EPI distributions versus task accuracies. β is slope coefficient of linear regression, r is correlation, and p is the two-tailed p value.

a standard isotropic gaussian $z_0 \sim \mathcal{N}(\mathbf{0}, I)$ mapped to a support of $\mathbf{z}_i \in [-5, 5]$. We used an augmented Lagrangian coefficient of $c_0 = 10^2$, a batch size $n = 100$, set $\nu = 0.5$, and initialized $q_{\boldsymbol{\theta}}(\mathbf{z})$ to produce an isotropic gaussian with mean 0 and variance 2.5^2. Accuracies were estimated over 200 trials of random gaussian noise, which was sampled independently for each drawn parameter $\mathbf{z}$ and each iteration of the EPI optimization.

5.2.4 Rank-2 RNN

Traditional approaches to likelihood-free inference – approximate Bayesian computation (ABC) methods – randomly sample parameters $\mathbf{z}$ until a suitable set is obtained. State-of-the-art ABC methods leverage sequential Monte Carlo (SMC) sampling techniques to obtain parameter sets more efficiently. To obtain more parameter samples, SMC-ABC must be run from scratch again. ABC methods do not confer log probabilities of samples. Like EPI, sequential neural posterior estimation (SNPE) uses deep learning to produce flexible posterior approximations. Like traditional Bayesian inference methods, SNPE conditions directly on the statistics of data. This differs from EPI, where posteriors are conditioned on emergent properties (moment constraints on the posterior predictive distribution). Peculiarities of SNPE (density estimation approach, two deep networks) make scaling

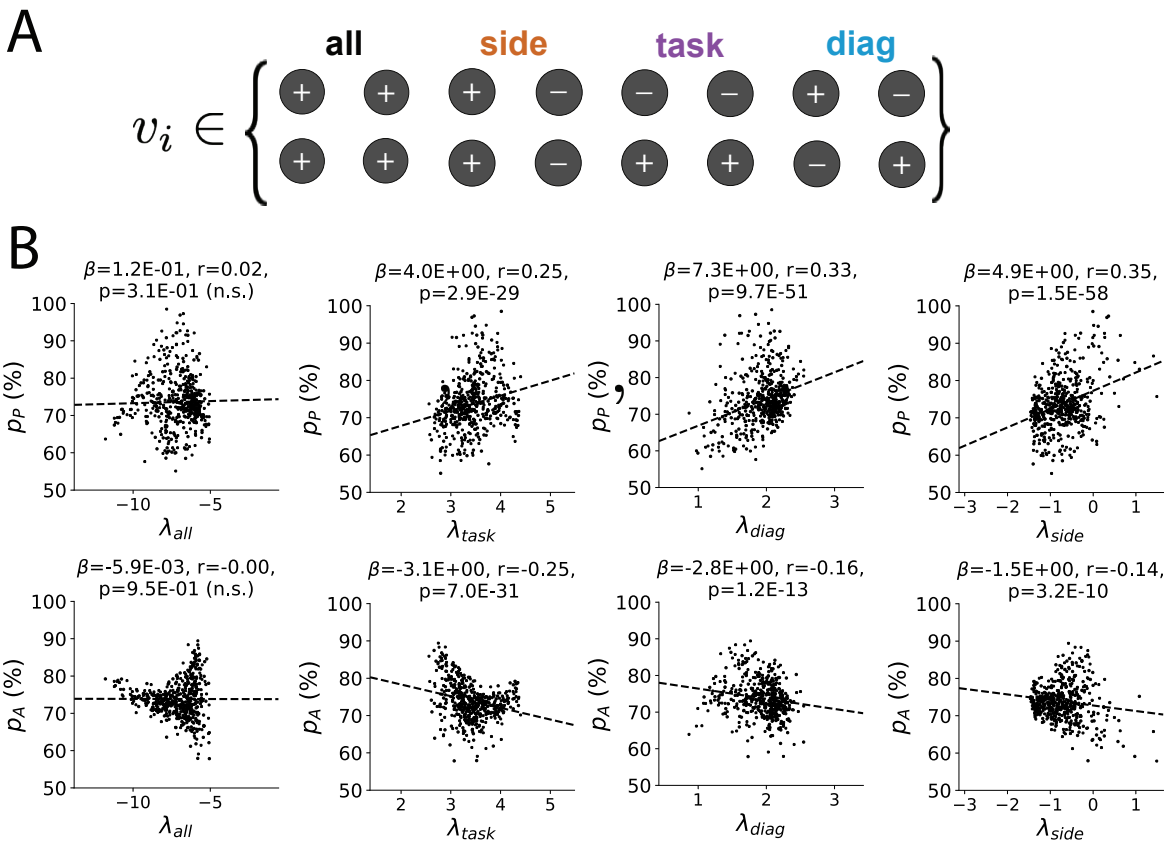

Figure 12: (SC2): A. Invariant eigenvectors of connectivity matrix W. B. Eigenvalues of connectivities of EPI distribution versus task accuracies.

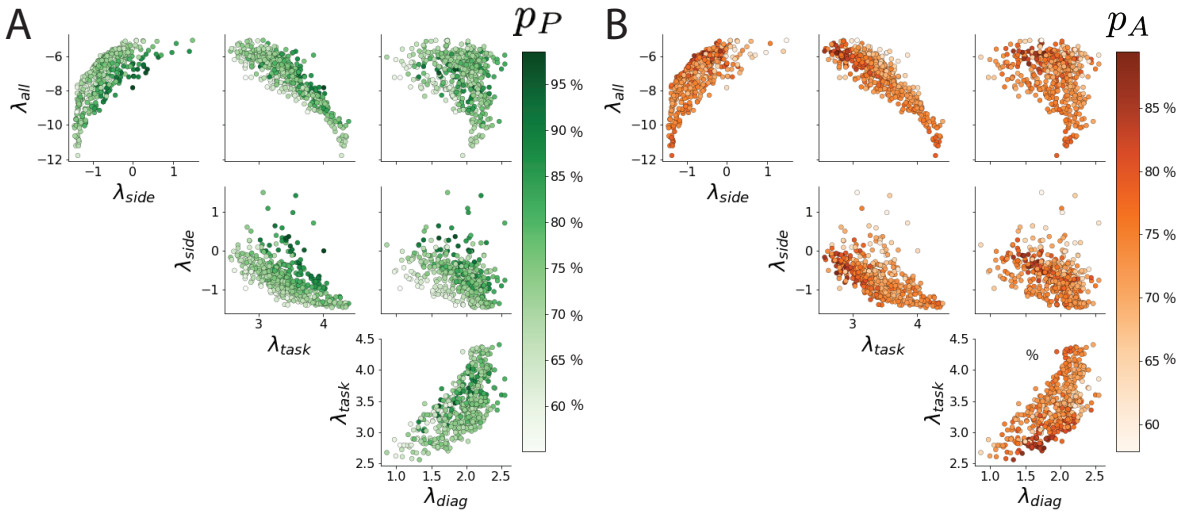

Figure 13: (SC3): A. Connectivity eigenvalues of EPI parameter distribution colored by Pro task accuracy. B. Same for Anti task.

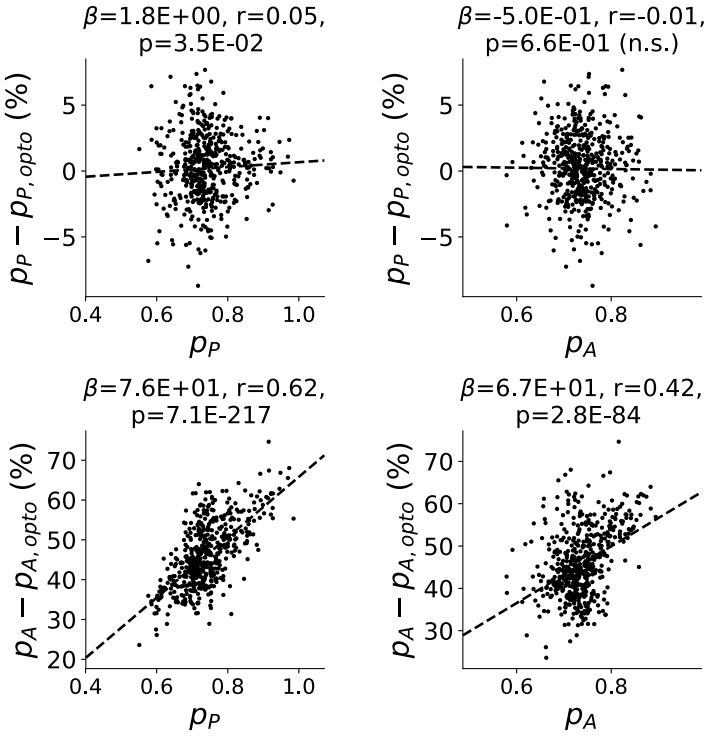

Figure 14: (SC4): Scatters of the effect of delay period inactivation in each task with task accuracy. Plots are shown at an opto strength of 0.8.

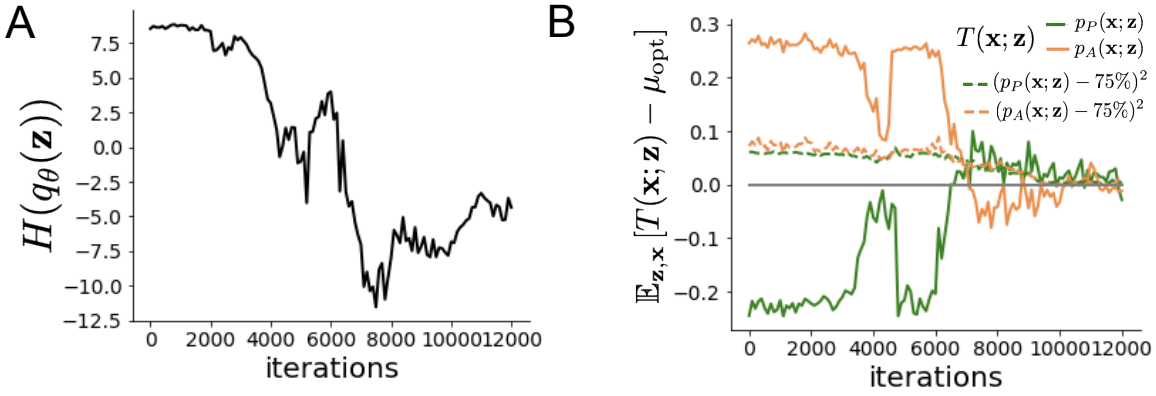

Figure 15: (SC5): EPI optimization of the SC model producing rapid task switching. A. Entropy throughout optimization. B. The emergent property statistic means and variances converge to their constraints at 12,000 iterations following the sixth augmented Lagrangian epoch.

in $\mathbf{z}$ prohibitive.

SMC-ABC has many hyperparameters, of which pyABC selects automatically by running some initial diagnostics upon initialization. In concurrence with the literature, SMC-ABC fails to converge around 25-30 dimensions, since it's proposal samples never get close enough to the target statistics. We searched over many SNPE hyperparameter choices: $n_{\text{train}} \in [2,000, 10,000, 100,000]$ is the number of simulations run per training epoch, and $n_{\text{mades}} \in [2, 3]$ is the number of masked autoregressive density estimators in the deep parameter distribution architecture. The greater n_{train}, the longer each epoch will take, but the more likely SNPE may converge during that epoch. Greater n_{mades} increases the flexibility of the deep parameter distribution of SNPE, but slows optimization. For the timing plot, we show the fastest among all of these choices, and for the convergence plot, we show the best convergence among all of these choices. During optimization, we used n_{atom}=100 atomic proposals as is recommended.