
Non-degenerate Priors for Arbitrarily Deep Networks

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Abstract

Choosing appropriate architectures and initialization strategies are crucial to good performance of deep networks. To shed light on this problem, we analyze the analogous problem of constructing useful priors on deep functions. We show that for typical deep architectures, the representational capacity of the network tends to capture fewer degrees of freedom as the number of layers increases, retaining only a single degree of freedom in the limit. We propose alternate priors on network structure which do not suffer from these pathologies, and also derive novel kernels obtained by composing infinitely-many feature transforms.

1 Introduction

Deep Gaussian processes are a prior on functions of the form: $f_{deep} = f^{1:L}(\mathbf{x}) = f^L(f^{L-1}(\dots f^3(f^2(f^1(\mathbf{x}))) \dots))$, where each $f \stackrel{\text{iid}}{\sim} \text{GP}$.

2 Relation between deep nets and Gaussian processes

Neal [1995] showed that the limit of a neural network with infinitely many hidden units. More precisely, he showed that for models of the form

$$f(x) = \alpha^T \Phi(x) = \sum_{i=1}^K \alpha_i \phi_i(x), \quad (1)$$

the central limit theorem implies that any two function values $f(x), f(x')$ have a distribution approaching $\mathcal{N}(\cdot, \cdot)$ as $K \rightarrow \infty$ with fixed feature vector $\Phi(x)$ and i.i.d. α approaches a distribution over functions in which all $f(x), f(x')$ are jointly Gaussian distributed, under broad conditions. The result is surprisingly general: It doesn't put any constraints on what the features are (other than having bounded activation), nor does it require that the feature weights α be Gaussian distributed. It only requires that the distribution on weights α has finite variance, and that as we increase the number of features K , we divide the feature weights by K . As $K \rightarrow \infty$, the central limit theorem gives that the weighted sum of feature activations approaches a Gaussian distribution.

2.1 A neural net with one hidden layer

In a neural net with a single hidden layer, the correspondence is simple: The features $\phi(\mathbf{x})$ correspond to the hidden units. In the typical definition of a multi-layer perceptron, the hidden units of the first layer are defined as:

$$\Phi(\mathbf{x}) = h^{(1)}(\mathbf{x}) = \sigma(b^{(1)} + W^{(1)}\mathbf{x}) \quad (2)$$

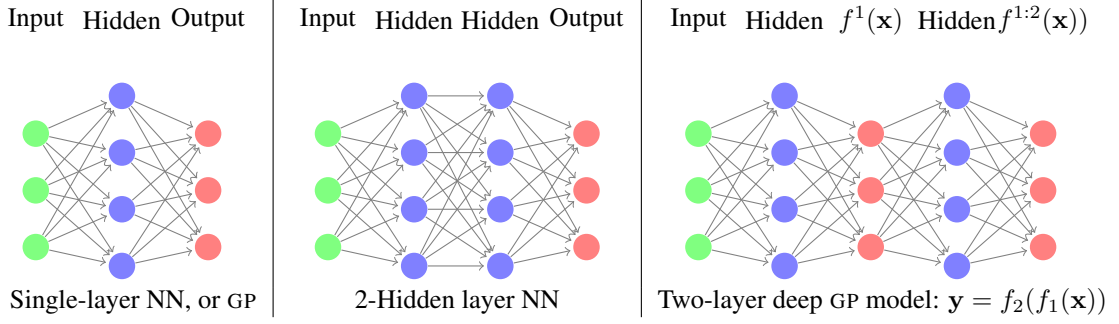


Figure 1: Comparing architectures. When there are multiple layers, the meaning of “hidden units” is ambiguous in the deep GP model. We can either consider every second layer to be a linear combination of an infinite number of parametric hidden basis functions, or we can ignore the hidden layers and consider the deep GP to be a neural network with a finite number of hidden units, with non-parametric activation functions.

The output $f(\mathbf{x})$ is simply a weighted sum of these hidden unit activations, for both GP models and neural networks.

2.2 A neural net with two hidden layers

In a neural net with two hidden layers, the correspondence is a little more complicated. Generally, the n^{th} layer units are given by the recurrence:

$$h^{(n)}(\mathbf{x}) = \sigma(b^{(n)} + W^{(n)}h^{(n-1)}(\mathbf{x})) \quad (3)$$

for real-valued outputs, the function can be written as weighted sum of the last layer’s hidden units:

$$f(\mathbf{x}) = W^{(top)}\sigma(b^{(n)} + W^{(n)}h^{(n-1)}(\mathbf{x})) = W^{(top)}h^{(n)}(\mathbf{x}) \quad (4)$$

We can also write deep GP models in this way. However, for many covariance functions, we can’t explicitly compute the activation of infinitely-many hidden functions $\Phi(\mathbf{x})$, and instead can only examine their weighted sum. Hence, to be able to compare the two model classes, we will view deep GP models in a different way, as a neural network with a finite number of nonparametric basis functions:

$$f^{1:L}(\mathbf{x}) = f^L(f^{L-1}(\dots f^3(f^2(f^1(\mathbf{x})))\dots)) \quad (5)$$

This view lets us compare GP models to multilayer perceptrons more directly, examining the activations and shapes of the finite number of basis functions at each layer, denoted by the vector of functions $f^\ell(\mathbf{x})$. Figure 1 compares these two architectures.

3 The relation between initialization strategies, regularization, and priors

Traditional training of deep networks requires both an initialization strategy, and a regularization method. In a Bayesian model, the prior implicitly defines both of these things.

3.1 What kinds of functions do deep GPs prefer?

Isotropic kernels such as the squared-exp encode the assumption that the function varies independently in all directions. Thus, deep GPs prefer to have hidden layers on whose outputs the higher-layer functions vary independently. In this way, we can say that the deep GP model prefers independent features.

The effects of automatic relevance determination One feature of Gaussian process models that doesn’t have a clear analogue in the neural-net literature is automatic relevance determination (ARD). Recall that in the ARD procedure, the lengthscales of each dimension are scaled to maximize the marginal likelihood of the GP. In standard one-layer GP regression, it has the effect of smoothing out the posterior over functions in directions in which the data does not change.

4 The Jacobian of Deep GPs is a product of i.i.d. Normal Matrices

In this section, we show a result which will let us characterize the Jacobians of arbitrarily deep networks drawn from a deep GP prior.

Lemma 4.1. *The partial derivatives of a function drawn from a GP prior with a product kernel are i.i.d. Gaussian distributed.*

Proof. Because differentiation is a linear operator, the derivatives of a function drawn from a GP prior are also jointly Gaussian distributed. The covariance between partial derivatives w.r.t. input dimensions d_1 and d_2 of vector \mathbf{x} are given by [Solak et al., 2003]:

$$\text{cov} \left(\frac{\partial f(\mathbf{x})}{\partial x_{d_1}}, \frac{\partial f(\mathbf{x})}{\partial x_{d_2}} \right) = \frac{\partial^2 k(\mathbf{x}, \mathbf{x}')}{\partial x_{d_1} \partial x'_{d_2}} \Big|_{\mathbf{x}=\mathbf{x}'} \quad (6)$$

If our kernel is a product over individual dimensions $k(\mathbf{x}, \mathbf{x}') = \prod_d k_d(x_d, x'_d)$, as in the case of the isotropic squared-exp kernel, then the off-diagonal entries are zero, implying that all elements are independent and identically distributed. \square

In the case of the isotropic squared-exp kernel, the covariance between derivatives has the form:

$$k_{SE}(\mathbf{x}, \mathbf{x}') = \sigma_f^2 \prod_{d=1}^D \left(-\frac{1}{2} \frac{(x_d - x'_d)^2}{\ell_d^2} \right) \implies \text{cov} \left(\frac{\partial f(\mathbf{x})}{\partial x_{d_1}}, \frac{\partial f(\mathbf{x})}{\partial x_{d_2}} \right) = \begin{cases} \frac{\sigma_f^2}{\ell_{d_1}^2} & \text{if } d_1 = d_2 \\ 0 & \text{if } d_1 \neq d_2 \end{cases} \quad (7)$$

Lemma 4.2. *The Jacobian of a GP with a seperable kernel is a matrix of i.i.d. Gaussian R.V.'s*

Proof. The Jacobian of the vector-valued function $f(\mathbf{x})$ is a matrix J whose i, j^{th} element is $J_{ij} = \frac{\partial f_i(\mathbf{x})}{\partial x_j}$. When composing L functions, we'll denote the *immediate* Jacobian of the function mapping from layer $\ell - 1$ to layer ℓ as $J^\ell(\mathbf{x})$, and the Jacobian of the entire composition of L functions by $J^{1:L}(\mathbf{x})$.

Because we've assumed that the GP on each output dimension $f_d(\mathbf{x}) \sim \mathcal{GP}$ is independent, it follows that for a given \mathbf{x} , each row of $J_{\mathbf{x} \rightarrow \mathbf{y}}(\mathbf{x})$ is independent. Above, we showed that the elements of each row are independent. This means that each entry in the Jacobian of a GP-distributed transform is i.i.d. Normal. \square

Theorem 4.3. *The Jacobian of a deep GP using a product kernel is a product of independent random Gaussian matrices.*

Proof. By the multivariate chain rule, the Jacobian of a composition of functions is simply the product of the Jacobian matrices of each function. Thus the Jacobian of the composed (deep) function $f^L(f^{L-1}(\dots f^3(f^2(f^1(\mathbf{x}))) \dots))$ is:

$$J^{1:L}(x) = J^L J^{L-1} \dots J^3 J^2 J^1 \quad (8)$$

Since each $J^i \stackrel{\text{iid}}{\sim} \mathcal{N}(\cdot)$, the complete Jacobian is a product of i.i.d. Gaussian matrices. \square

This result means that we can analyze the representational properties of a deep Gaussian process by simply examining the properties of products of i.i.d. Gaussian matrices.

5 Formalizing the pathologies

5.1 Desirable properties of latent representations

Rifai et al. [2011a] make the point that a good latent representation is one that remains invariant when moving in directions orthogonal to the manifold that the data lie on. Conversely, a good latent representation must also change in directions tangential to the data manifold - otherwise we are losing information about the data. Figure 2 demonstrates this idea.

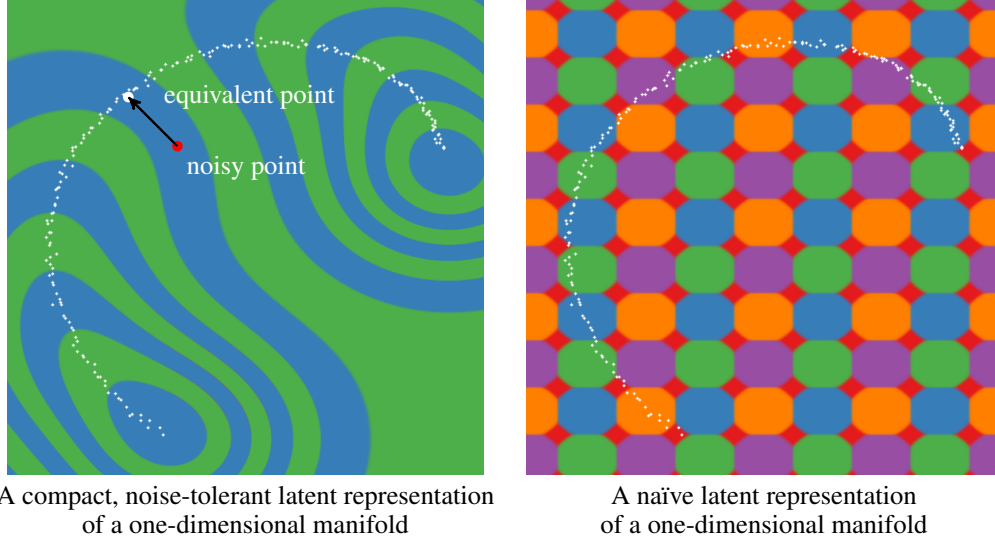


Figure 2: Comparing different latent representations of data on a 1-D manifold. A good latent representation is invariant in directions orthogonal to the data manifold, making the representation robust to noise in those directions. The representation must also change in directions tangent to the manifold, in order to preserve information for later layers.

We follow Rifai et al. [2011b] in characterizing the representational properties of a function by the singular value spectrum of the Jacobian¹.

One-dimensional Asymptotics For a one-dimensional deep function, the derivative is simply a product of i.i.d Normal variables, with zero mean and variance $\frac{\sigma_f^2}{\ell^2}$. The distribution of the absolute value of the derivative of the deep function is a product of half-normals:

$$\frac{\partial f(x)}{\partial x} \stackrel{iid}{\sim} \mathcal{N}\left(0, \frac{\sigma_f^2}{\ell^2}\right) \implies \left|\frac{\partial f(x)}{\partial x}\right| \stackrel{iid}{\sim} \text{half}\mathcal{N}\left(\sqrt{\frac{2\sigma_f^2}{\pi\ell^2}}, \frac{\sigma_f^2}{\ell^2} \left(1 - \frac{2}{\pi}\right)\right) \quad (9)$$

Thus if $\frac{\sigma_f^2}{\ell^2} = \frac{\pi}{2}$, then $\mathbb{E}\left[\left|\frac{\partial f(x)}{\partial x}\right|\right] = 1$, and so $\mathbb{E}\left[\left|\frac{\partial f^{1:L}(x)}{\partial x}\right|\right] = 1$. If $\frac{\sigma_f^2}{\ell^2}$ is less than $\frac{\pi}{2}$, the expected derivative magnitude goes to zero, and if it's greater, then the expected magnitude goes to infinity.

The log of this variable has moments:

$$m_{\log} = \mathbb{E}\left[\log\left|\frac{\partial f(x)}{\partial x}\right|\right] = 2\log\left(\frac{\sigma_f}{\ell}\right) - \log 2 - \gamma \quad (10)$$

$$v_{\log} = \mathbb{V}\left[\log\left|\frac{\partial f(x)}{\partial x}\right|\right] = \frac{\pi^2}{4} + \frac{\log^2 2}{2} - \gamma^2 - \gamma \log 4 + 2\log\left(\frac{\sigma_f}{\ell}\right) \left[\gamma + \log 2 - \log\left(\frac{\sigma_f}{\ell}\right)\right] \quad (11)$$

Since the second moment is finite, by the central limit theorem, the limiting distribution of the size of the gradient is log-normal:

$$\log\left|\frac{\partial f^{1:L}(x)}{\partial x}\right| = \sum_{i=1}^L \log\left|\frac{\partial f^L(x)}{\partial x}\right| \quad (12)$$

$$\log\left|\frac{\partial f^{1:L}(x)}{\partial x}\right| \stackrel{L \rightarrow \infty}{\sim} \mathcal{N}(Lm_{\log}, L^2v_{\log}) \quad (13)$$

¹Rifai et al. [2011b] examine the Jacobian at the training points, but the models we are examining are stationary, so it doesn't matter where we examine the function.

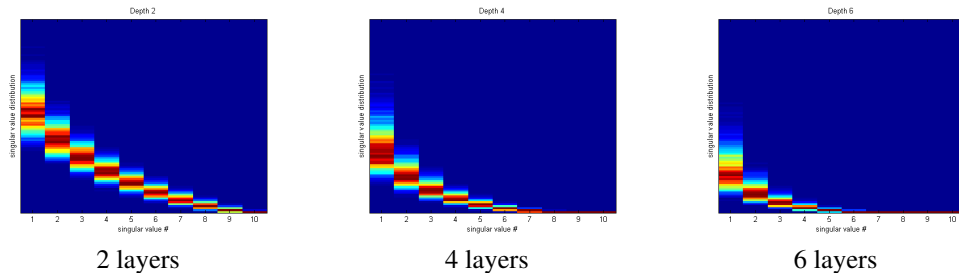


Figure 3: Singular value spectrum of the Jacobian of a deep GP. As the net gets deeper, the largest singular value dominates. This implies that there is only one effective degree of freedom in representation being computed.

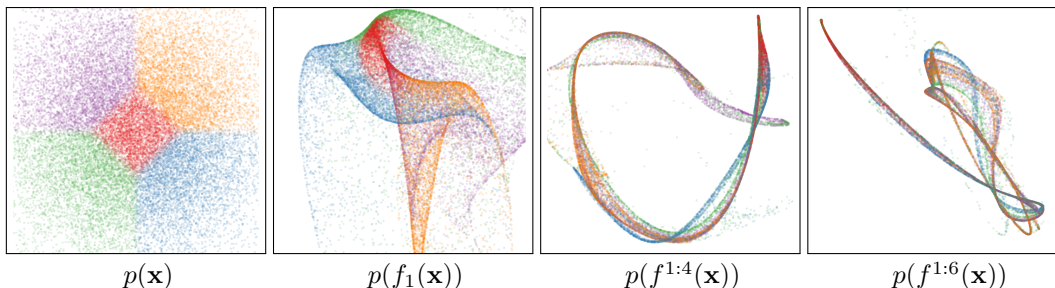


Figure 4: Draws from a deep GP. A distribution is warped by successive functions drawn from a GP prior. As the number of layers increases, the density exhibits a sort of filamentation.

Even if the expected magnitude of the derivative remains constant, the variance of the distribution grows without bound, and so will almost surely be very small almost everywhere, with rare but very large jumps. Note that there is another limit - if $\frac{\sigma_f^2}{\ell_{d_1}^2} < \frac{\pi}{2}$, then the mean of the size of the gradient goes to zero, but the variance approaches a finite constant.

With high probability, the regions of $f^{1:L}(\mathbf{x})$ which vary quickly will be the same regions in which $f^{1:L+1}(\mathbf{x})$ has high variance.

5.2 Jacobian Singular Value Spectrum

Figure 3 shows the spectrum for deep GPs of different depths. As the net gets deeper, the largest singular value dominates. This implies that there is only one effective degree of freedom in representation being computed. As well, the distribution on singular values seems to become heavy-tailed.

Figure 4 demonstrates a related pathology that arises when composing functions to produce a deep latent-variable model. The density in observed space eventually becomes locally concentrated onto one-dimensional manifolds, or *filaments*.

To visualize this pathology in another way, figure 5 illustrates the value computed at each point in the input space, after successive warpings. After 40 warpings, we can see that locally, there is usually only one direction that one can move in \mathbf{x} -space in order to change the value of the function.

5.3 Fixing the pathology

Follow a suggestion in Neal [1995], we can fix the pathologies exhibited in figures 4 and 5 by simply making each layer of computation depend not only on the output of the previous layer, but also on the original input \mathbf{x} . Draws from the resulting priors are shown in figures 8 and 9.

Figure 6 shows draws from a 1D deep prior, both with naive compositions, and connected compositions.

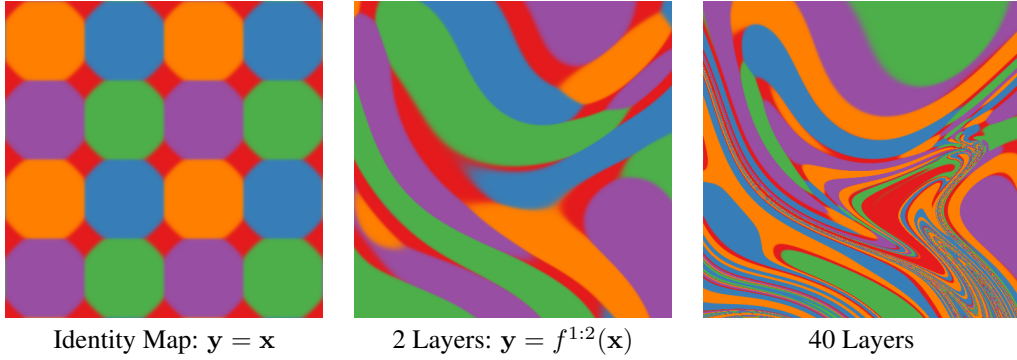


Figure 5: Feature Mapping of a deep GP. Shown here are the colors corresponding to the location $y = f(x)$ that each point is mapped to after being warped by a deep GP. Just as the densities in 4 became locally one-dimensional, there is locally only one direction that one can move x in to change y . Regions where the colors change rapidly indicate that the Jacobian is large in that area.

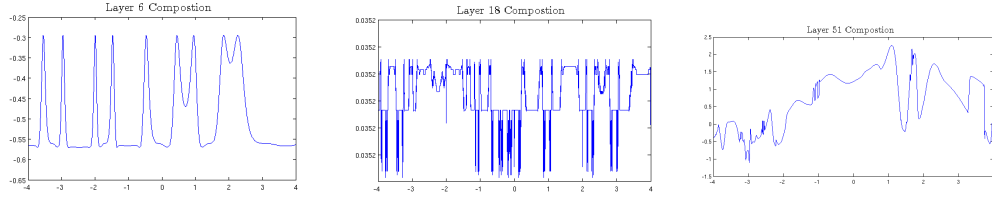


Figure 6: One-dimensional draws from deep Gaussian processes. The two left draws are from the naive compositional architecture, and rapidly become degenerate. The right-hand draw has every layer connected to the input, and maintains smoothness in some regions, while varying rapidly in other regions.

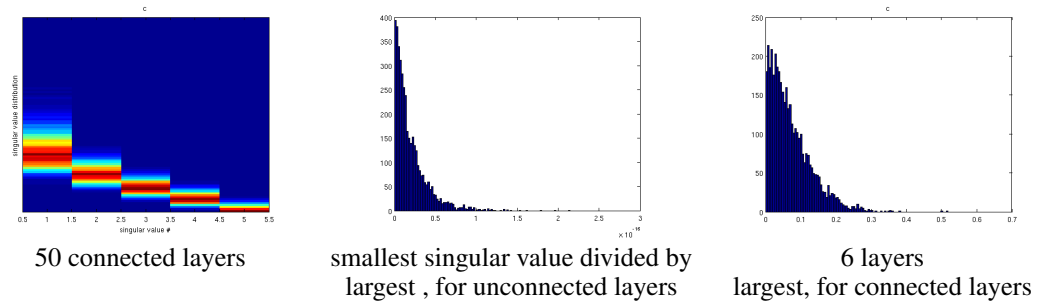


Figure 7: Repairing the singular value distribution. When each layer is connected to the original inputs, and the variance of derivatives is smaller than a threshold, then the ratio of the largest singular value to the smallest remains small.

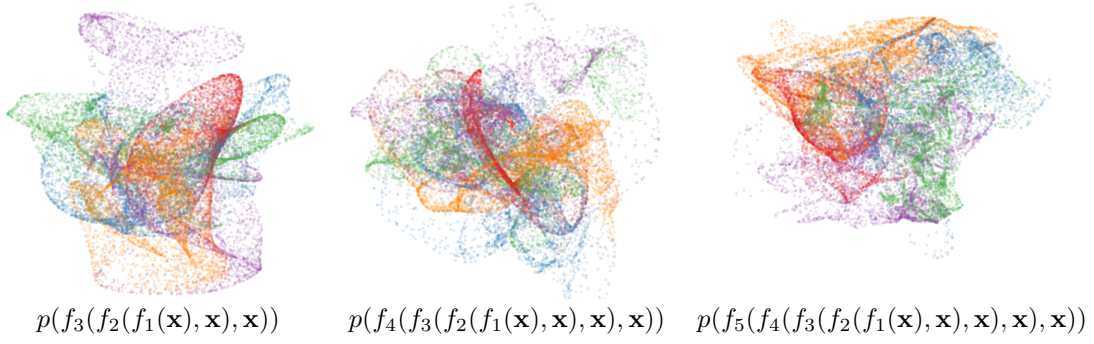


Figure 8: Draws from a deep GP, with each layer connected to the input \mathbf{x} . By always depending on the original input, the density becomes more complex without concentrating along filaments.

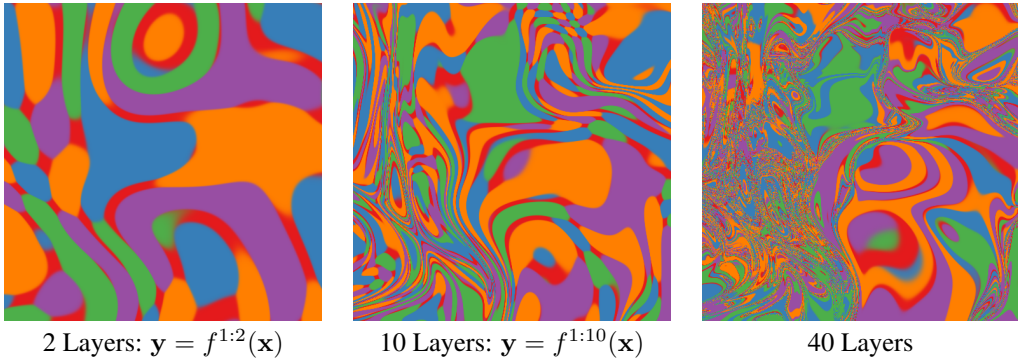


Figure 9: Feature Mapping of a deep GP with each layer connected to the input \mathbf{x} . Just as the densities in 8 become remain locally two-dimensional even after many transformations, in this mapping there are locally usually two directions that one can move \mathbf{x} in to change \mathbf{y} .

Jacobians of connected deep networks We can similarly examine the Jacobians of the new connected architecture. The Jacobian of a composite function which has had the original inputs appended $f_{\text{aug}}(\mathbf{x}) = [f(\mathbf{x}), \mathbf{x}]$, is $\begin{bmatrix} J_f \\ I_D \end{bmatrix}$ and the Jacobian of all one-layer transformations of these augmented functions are $D \times 2D$ matrices. The Jacobian of the composed, connected deep function is defined by the recurrence:

$$J^{1:L}(\mathbf{x}) = J^L \begin{bmatrix} J^{1:L-1} \\ I_D \end{bmatrix} \quad (14)$$

5.4 Sparse Initialization

Martens [2010] used sparse initialization, in which each hidden unit had 15 non-zero incoming connections. They said that “Doing this allows the units to be both highly differentiated as well as unsaturated, avoiding the problem in dense initializations where the connection weights must all be scaled very small in order to prevent saturation, leading to poor differentiation between units”.

While it is not clear if deep GPs have analogous problems with saturation of connection weights, we note that an analogous initialization strategy would be to put a heavy-tailed prior on the lengthscales in each kernel.

6 Arbitrarily Deep Kernels

Cho [2012] investigated kernels constructed by applying multiple layers of feature mappings. That is to say, if a kernel has the form $k_1(\mathbf{x}, \mathbf{x}') = \Phi(\mathbf{x})^\top \Phi(\mathbf{x}')$, then we can consider constructing a kernel based on the repeated feature mapping: $k_2(\mathbf{x}, \mathbf{x}') = \Phi(\Phi(\mathbf{x}))^\top \Phi(\Phi(\mathbf{x}'))$.

This composition operation has a closed form for any set of starting features $\Phi_n(\mathbf{x})$:

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp \left(-\frac{1}{2} \|\Phi_n(\mathbf{x}) - \Phi_n(\mathbf{x}')\|_2^2 \right) \quad (15)$$

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp \left(-\frac{1}{2} \sum_i \left[\phi_n^{(i)}(\mathbf{x}) - \phi_n^{(i)}(\mathbf{x}') \right]^2 \right) \quad (16)$$

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp \left(-\frac{1}{2} [k_n(\mathbf{x}, \mathbf{x}) - 2k_n(\mathbf{x}, \mathbf{x}') + k_n(\mathbf{x}', \mathbf{x}')] \right) \quad (17)$$

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp(k_n(\mathbf{x}, \mathbf{x}') - 1) \quad (\text{if } k_n(\mathbf{x}, \mathbf{x}) = 1) \quad (18)$$

Note that nothing in this derivation depends on details of k_n , except that $k_n(\mathbf{x}, \mathbf{x}) = 1$.

6.1 Infinitely deep kernels

What happens when we apply this composition many times, starting with the squared-exp kernel? In the infinite limit, this recursion converges to $k(x, y) = 1$ for all inputs. Figure ?? shows this kernel at different depths, including the degenerate limit.

One interpretation of why repeated feature transforms lead to this degenerate prior is that each layer can only lose information about the previous set of features. In the limit, the transformed features contain no information about the original input \mathbf{x} . Since the function doesn't depend on its input, it must be the same everywhere.

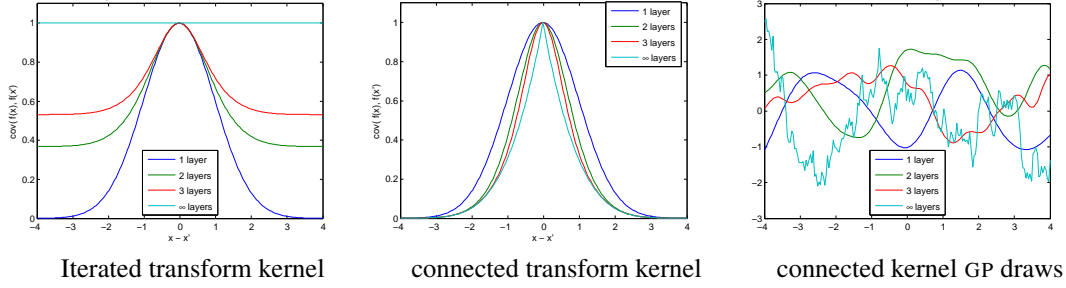


Figure 10: A non-degenerate version of the infinitely deep feature transform kernel. By connecting the inputs \mathbf{x} to each layer, the function can still depend on its input even after arbitrarily many layers of computation.

6.2 A non-degenerate construction

Following a suggestion from Neal [1995], we append the inputs \mathbf{x} to each layer of features. To do so, we simply by augmenting the feature vector $\Phi_n(\mathbf{x})$ with the extra features \mathbf{x} at each layer:

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp \left(-\frac{1}{2} \left\| \begin{bmatrix} \Phi_n(\mathbf{x}) \\ \mathbf{x} \end{bmatrix} - \begin{bmatrix} \Phi_n(\mathbf{x}') \\ \mathbf{x}' \end{bmatrix} \right\|_2^2 \right) \quad (19)$$

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp \left(-\frac{1}{2} \sum_i [\phi_i(\mathbf{x}) - \phi_i(\mathbf{x}')]^2 - \frac{1}{2} \|\mathbf{x} - \mathbf{x}'\|_2^2 \right) \quad (20)$$

$$k_{n+1}(\mathbf{x}, \mathbf{x}') = \exp \left(k_n(\mathbf{x}, \mathbf{x}') - 1 - \frac{1}{2} \|\mathbf{x} - \mathbf{x}'\|_2^2 \right) \quad (21)$$

This kernel satisfies the recurrence $k - \log(k) = 1 + \frac{1}{2} \|\mathbf{x} - \mathbf{x}'\|_2^2$, a non-degenerate limit. The solution to this recurrence has no closed form, but it is continuous and differentiable everywhere except at $\mathbf{x} = \mathbf{x}'$. Samples from a GP with this prior are not differentiable, and locally resemble brownian motion, having a sort of fractal behavior.

6.3 Can “Deep Kernels” ever be powerful models?

Bengio et al. [2006] showed that kernel machines, such as GPs, have limited generalization ability when they use a ‘local’ kernel such as the squared-exp. However, many interesting non-local kernels can be constructed which allow non-trivial extrapolation, for example, periodic kernels. A periodic kernel can be viewed as a 2-layer “deep kernel” [cite Mackay?], in which the first layer maps the input $x \rightarrow [\sin(x), \cos(x)]$, and the second layer maps through a set of radial-basis functions.

Other methods for constructing kernels can capture many other types of structure Duvenaud et al. [2013], Wilson and Adams [2013], such as translation and rotation invariance in images Kondor [2008]. Salakhutdinov and Hinton [2008] even uses a deep neural network to construct feature transforms for kernels.

6.4 More general feature compositions

The failure of the kernels derived in this section to produce interesting computations suggest that deep computation is only powerful when it is combined with learning.

However, we now know how to map any kernel’s feature mapping through an RBF feature mapping in closed form, and also how to append any other set of features to such a kernel. Thus we can now define kernels corresponding to arbitrary fixed computation graphs. This might be a nice way to extend kernel-search procedures.

7 Related Work

Variational inference in deep Gaussian processes was developed by [Damianou and Lawrence \[2012\]](#), who also analyzed the effect of ARD in these models.

Several variants of deep models have been proposed, such as Bayesian Deep Networks [Adams et al. \[2010\]](#) and Sum-product networks [Poon and Domingos \[2011\]](#). These models also have no connections except between adjacent layers, and perhaps may have similar pathologies as the number of layers increases.

Deep Density Networks [Rippel and Adams \[2013\]](#) were constructed with invertibility in mind, with parameters encouraged to preserve information about lower layers.

[Pascanu et al. \[2012\]](#) analyze the exploding-gradients problem in recurrent neural networks.

[Montavon et al. \[2010\]](#) note that performance of a MLP degrades as the number of layers with random weights increases.

8 Discussion

To what extent are these pathologies present in nets being used today? In simulations, we found that for deep functions with a fixed latent dimension D , that the singular value spectrum remained relatively flat for hundreds of layers as long as $D > 100$.

Recursive learning method Just as layer-wise unsupervised pre-training encourages the projection of the data into a representation with independent features in the higher layers, so does the procedure outlined here. This is because the isotropic kernel does not penalize independence between different dimensions, only the number of dimensions.

9 Conclusions

- Deep neural networks and deep Gaussian processes are analyzable using random matrix theory.
- If you want to use very deep nets, you won't be able to do so if you initialize/regularize all your weights independently.
- If you connect the input to every layer, some of these pathologies go away.

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