

# *Sparse Polynomial Chaos Surrogate for ACME Land Model*

SAND2015-10667C

## *via Iterative Bayesian Compressive Sensing*

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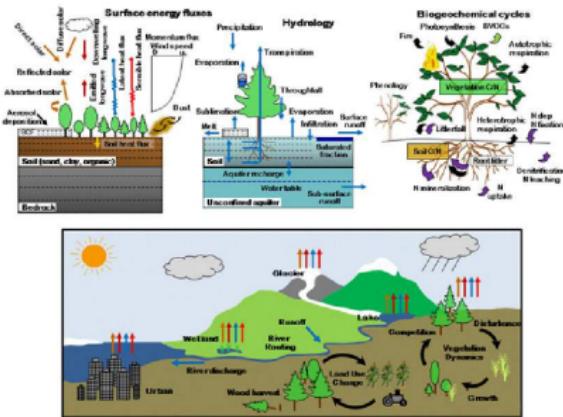
*Sponsored by DOE, Biological and Environmental Research,  
under Accelerated Climate Modeling for Energy (ACME).*

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# OUTLINE

- Surrogates needed for complex models
- Polynomial Chaos (PC) surrogates do well with uncertain inputs
- Bayesian regression provide results with uncertainty certificate
- Compressive sensing ideas deal with high-dimensionality

# Application of Interest: ACME Land Model



<http://www.cesm.ucar.edu/models/clm/>

- Nested computational grid hierarchy
- A single-site, 1000-yr simulation takes  $\sim 10$  hrs on 1 CPU
- Involves  $\sim 70$  input parameters; some dependent
- Non-smooth input-output relationship

# Surrogate construction: scope and challenges

Construct surrogate for a complex model  $f(\lambda)$  to enable

- Global sensitivity analysis
- Optimization
- Forward uncertainty propagation
- Input parameter calibration
- ...

- Computationally expensive model simulations, data sparsity
  - Need to build accurate surrogates with as few training runs as possible
- High-dimensional input space
  - Too many samples needed to cover the space
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# Polynomial Chaos surrogate for $f(\lambda)$

- Scale the input parameters  $\lambda_i \in [a_i, b_i]$

$$\lambda_i = \frac{a_i + b_i}{2} + \frac{b_i - a_i}{2} x_i$$

- Forward function  $f(\cdot)$ , output  $u$

$$u = f(\lambda(x)) \approx \sum_{k=0}^{K-1} c_k \Psi_k(x) \equiv g(x)$$

- Global sensitivity information for free
  - Sobol indices, variance-based decomposition.
- Bayesian inference useful for finding  $c_k$ :

$$P(c_k | u(x_j)) \propto P(u(x_j) | c_k) P(c_k)$$

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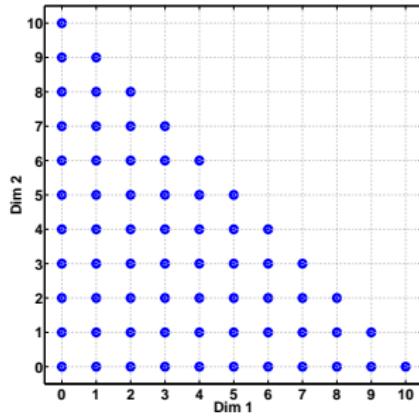
# Bayesian inference of PC surrogate: high-d, low-data regime

$$y = u(\mathbf{x}) \approx \sum_{k=0}^{K-1} c_k \Psi_k(\mathbf{x})$$

$$\Psi_k(x_1, x_2, \dots, x_d) = \psi_{k_1}(x_1) \psi_{k_2}(x_2) \cdots \psi_{k_d}(x_d)$$

- Issues:

- how to properly choose the basis set?
- need to work in underdetermined regime  $N < K$ : fewer data than bases (d.o.f.)

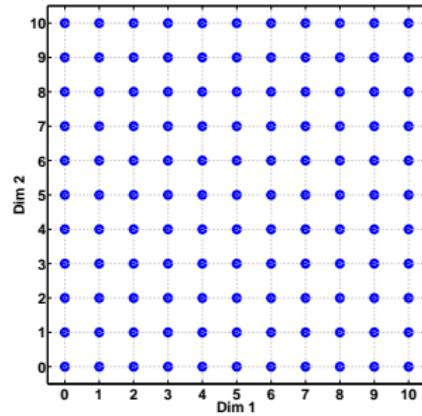


- Discover the underlying low-d structure in the model
  - get help from the machine learning community

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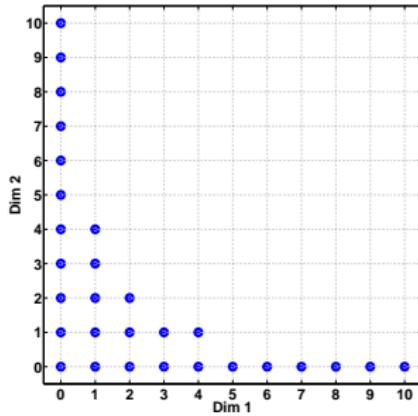
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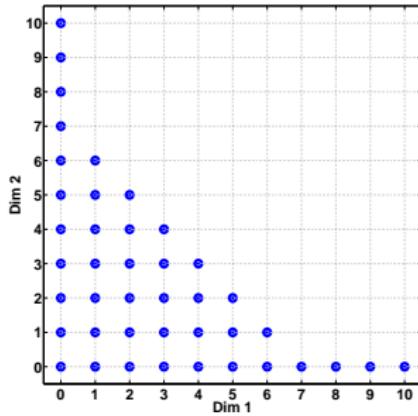


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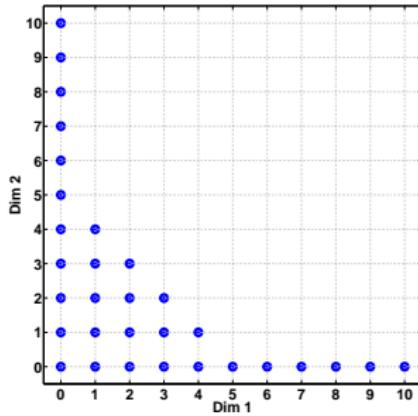


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# In a different language....

- $N$  training data points  $(\mathbf{x}_n, u_n)$  and  $K$  basis terms  $\Psi_k(\cdot)$
- Projection matrix  $\mathbf{P}^{N \times K}$  with  $\mathbf{P}_{nk} = \Psi_k(\mathbf{x}_n)$
- Find regression weights  $\mathbf{c} = (c_0, \dots, c_{K-1})$  so that

$$\mathbf{u} \approx \mathbf{P}\mathbf{c}$$

or

$$u_n \approx \sum_k c_k \Psi_k(\mathbf{x}_n)$$

- The number of polynomial basis terms grows fast; a  $p$ -th order,  $d$ -dimensional basis has a total of  $K = (p+d)!/(p!d!)$  terms.
- For limited data and large basis set ( $N < K$ ) this is a sparse signal recovery problem  $\Rightarrow$  need some regularization/constraints.

- Least-squares  $\text{argmin}_{\mathbf{c}} \{ \|\mathbf{u} - \mathbf{P}\mathbf{c}\|_2 \}$

- The ‘sparsest’  $\text{argmin}_{\mathbf{c}} \{ \|\mathbf{u} - \mathbf{P}\mathbf{c}\|_2 + \alpha \|\mathbf{c}\|_0 \}$

- Compressive sensing  $\text{argmin}_{\mathbf{c}} \{ \|\mathbf{u} - \mathbf{P}\mathbf{c}\|_2 + \alpha \|\mathbf{c}\|_1 \}$

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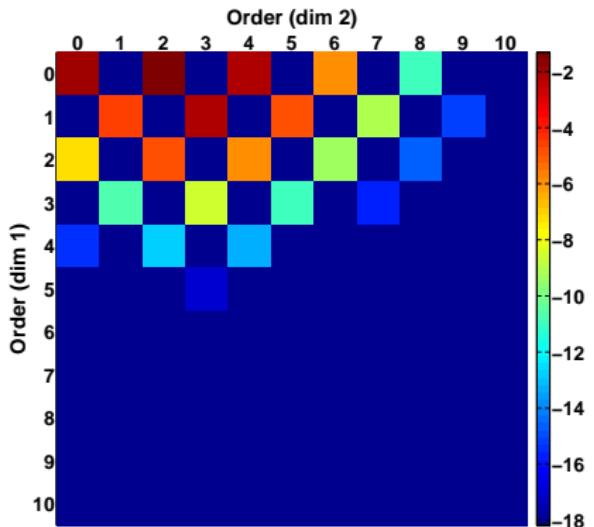
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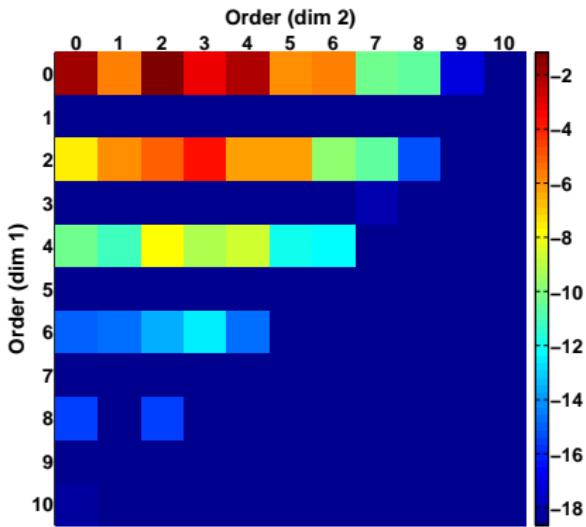
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Bayesian  $\text{argmin}_{\mathbf{c}} \{ \|\mathbf{u} - \mathbf{P}\mathbf{c}\|_2 + \alpha \|\mathbf{c}\|_1 \}$   
Likelihood Prior

# BCS removes unnecessary basis terms

$$f(x, y) = \cos(x + 4y)$$



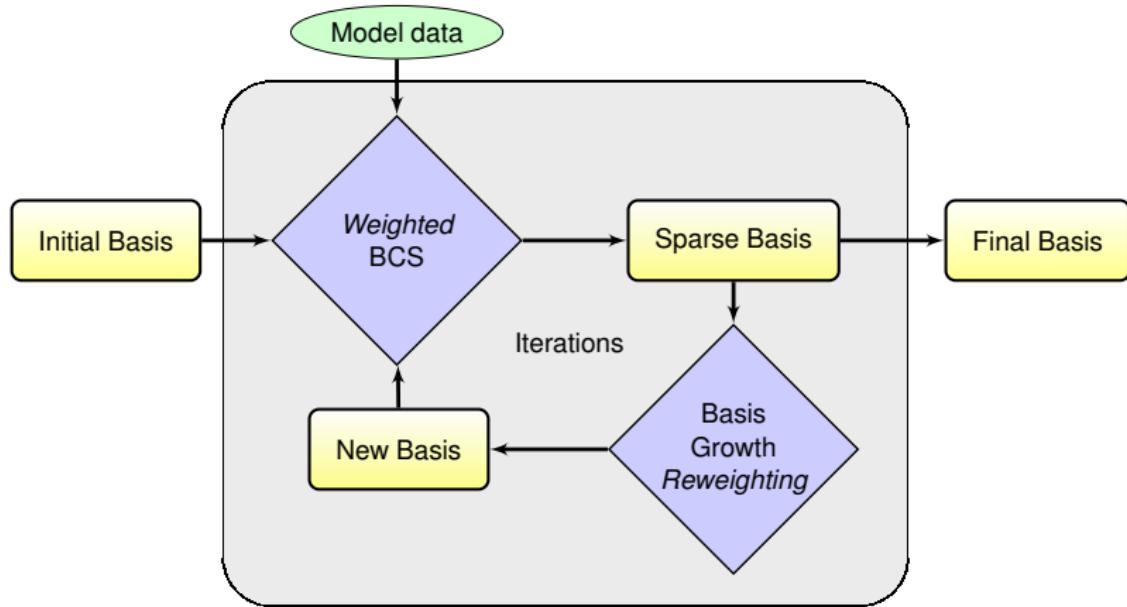
$$f(x, y) = \cos(x^2 + 4y)$$



The square  $(i, j)$  represents the (log) spectral coefficient for the basis term  $\psi_i(x)\psi_j(y)$ .

# Iterative Bayesian Compressive Sensing (iBCS)

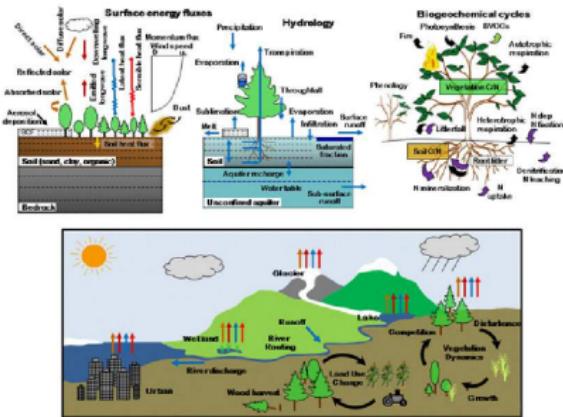
- *Iterative BCS*: We implement an iterative procedure that allows increasing the order for the relevant basis terms while maintaining the dimensionality reduction [Sargsyan *et al.* 2014], [Jakeman *et al.* 2015].
- Combine basis growth and reweighting!



# Basis set growth: simple anisotropic function

# Basis set growth: ... added outlier term

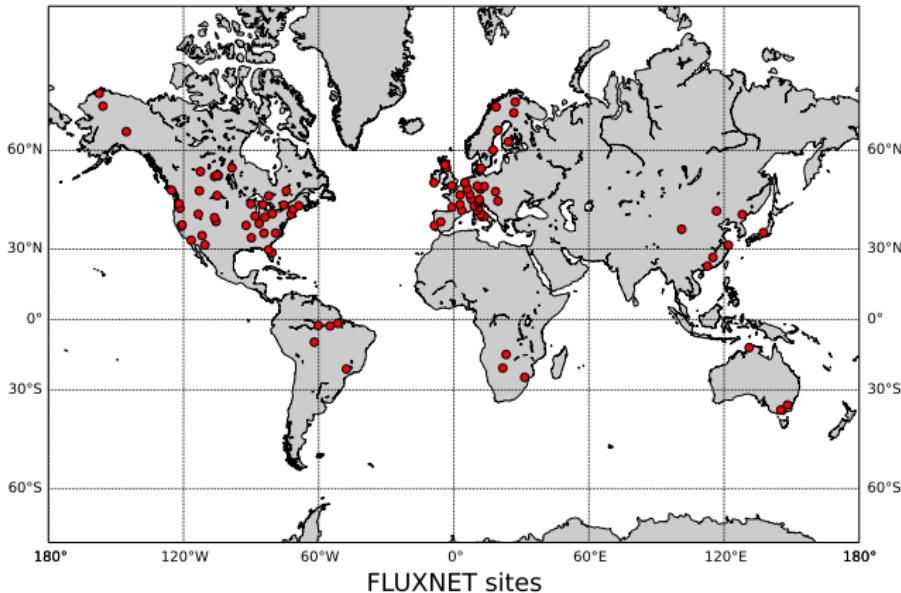
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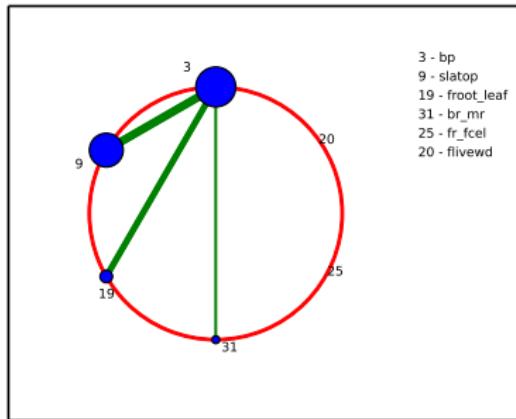
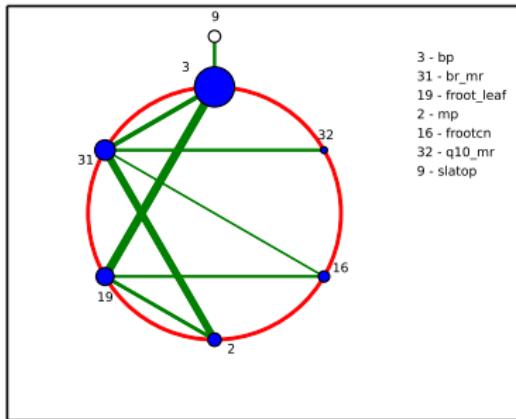
# FLUXNET experiment



- 96 FLUXNET sites covering major biomes and plant functional types
- Varying 68 input parameters over given ranges; 5 steady state outputs
- Ensemble of 3000 runs on Titan, DoE Leadership Computing Facility at Oak Ridge National Lab

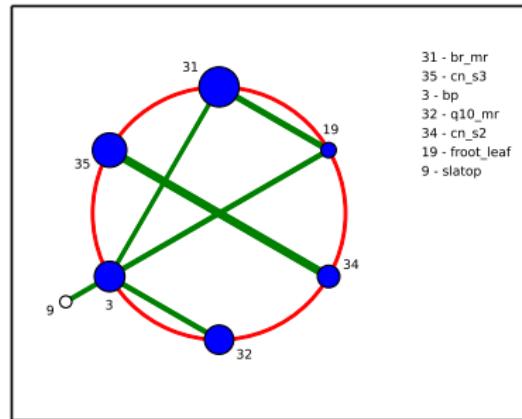
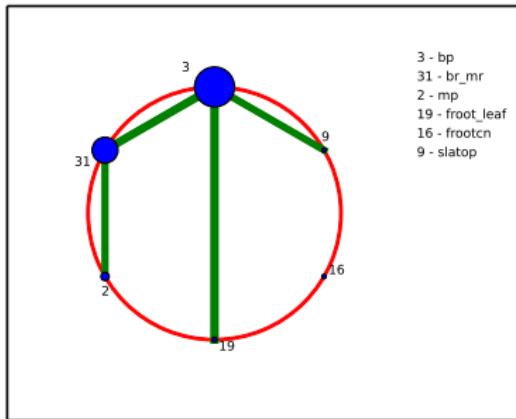
# Sparse PC surrogate and uncertainty decomposition for the ACME Land Model

- Main effect sensitivities : rank input parameters
- Joint sensitivities : most influential input couplings
- About 200 polynomial basis terms in the 68-dimensional space
- Sparse PC will further be used for
  - sampling in a reduced space
  - parameter calibration against experimental data



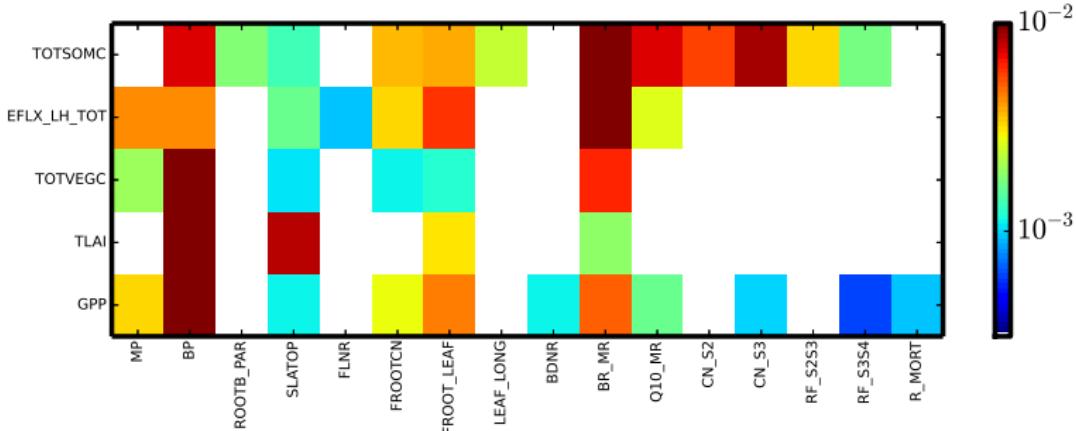
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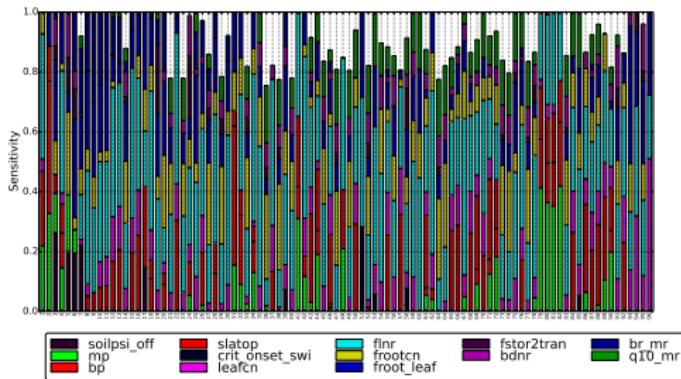
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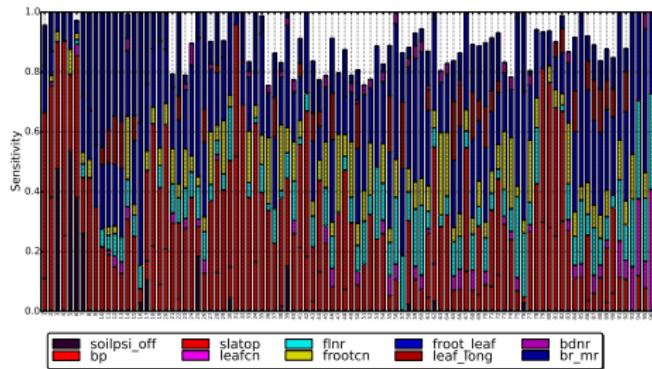
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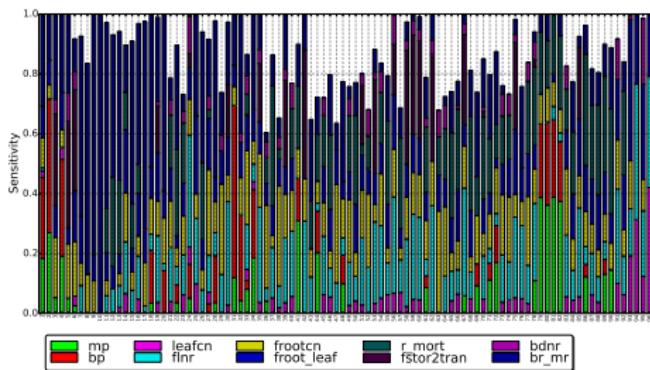
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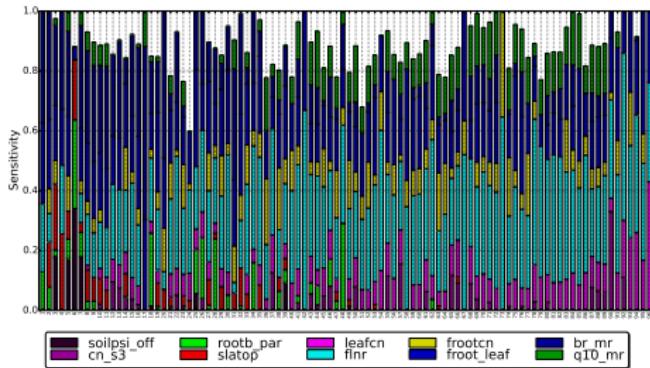
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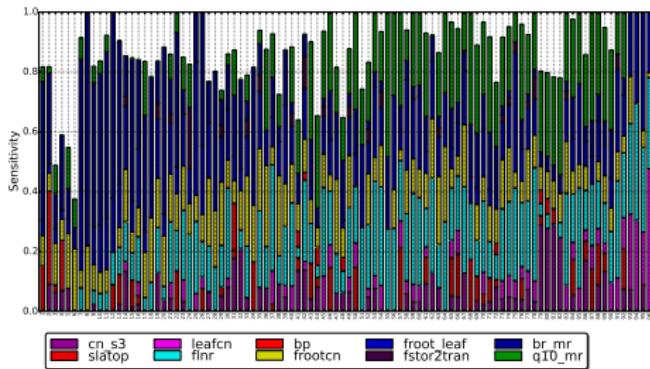
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# Summary

- **Surrogate** models are necessary for complex models
  - Replace the full model for both forward and inverse UQ
- Uncertain inputs
  - **Polynomial Chaos** surrogates well-suited
- Limited training dataset
  - **Bayesian** methods handle limited information well
- Curse of dimensionality
  - The hope is that not too many dimensions matter
  - Compressive sensing (CS) ideas ported from machine learning
  - We implemented *iteratively* **reweighting Bayesian CS** algorithm that reduces dimensionality and increases order on-the-fly.

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- Open issues
  - Computational design. What is the best sampling strategy?
  - Overfitting still present. Cross-validation techniques help.

# Literature

- M. Tipping, "Sparse Bayesian learning and the relevance vector machine", *J Machine Learning Research*, 1, pp. 211-244, 2001.
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- J. Jakeman, M. Eldred and K. Sargsyan, "Enhancing  $\ell_1$ -minimization estimates of polynomial chaos expansions using basis selection", *J Comp Phys*, 289, pp. 18-34, 2015.

# Random variables represented by Polynomial Chaos

$$X \simeq \sum_{k=0}^{K-1} c_k \Psi_k(\boldsymbol{\eta})$$

- $\boldsymbol{\eta} = (\eta_1, \dots, \eta_d)$  standard i.i.d. r.v.  
 $\Psi_k$  standard polynomials, orthogonal w.r.t.  $\pi(\boldsymbol{\eta})$ .

$$\Psi_k(\eta_1, \eta_2, \dots, \eta_d) = \psi_{k_1}(\eta_1) \psi_{k_2}(\eta_2) \cdots \psi_{k_d}(\eta_d)$$

- Typical truncation rule: total-order  $p$ ,  $k_1 + k_2 + \dots + k_d \leq p$ .  
Number of terms is  $K = \frac{(d+p)!}{d!p!}$ .
- Essentially, a parameterization of a r.v. by deterministic spectral modes  $c_k$ .
- Most common standard Polynomial-Variable pairs:  
(continuous) Gauss-Hermite, Legendre-Uniform,  
(discrete) Poisson-Charlier.

[Wiener, 1938; Ghanem & Spanos, 1991; Xiu & Karniadakis, 2002; Le Maître & Knio, 2010]

# Bayesian inference of PC surrogate

$$u \simeq \sum_{k=0}^{K-1} c_k \Psi_k(\mathbf{x}) \equiv g\mathbf{c}(\mathbf{x})$$

$$\overbrace{P(\mathbf{c}|\mathcal{D})}^{\text{Posterior}} \propto \overbrace{P(\mathcal{D}|\mathbf{c})}^{\text{Likelihood}} \overbrace{P(\mathbf{c})}^{\text{Prior}}$$

- Data consists of *training runs*

$$\mathcal{D} \equiv \{(\mathbf{x}_i, u_i)\}_{i=1}^N$$

- Likelihood with a gaussian noise model with  $\sigma^2$  fixed or inferred,

$$L(\mathbf{c}) = P(\mathcal{D}|\mathbf{c}) = \left( \frac{1}{\sigma \sqrt{2\pi}} \right)^N \prod_{i=1}^N \exp \left( -\frac{(u_i - g\mathbf{c}(\mathbf{x}))^2}{2\sigma^2} \right)$$

- Prior on  $\mathbf{c}$  is chosen to be conjugate, uniform or gaussian.
- Posterior is a *multivariate normal*

$$\mathbf{c} \in \mathcal{MVN}(\boldsymbol{\mu}, \boldsymbol{\Sigma})$$

- The (uncertain) surrogate is a *gaussian process*

$$\sum_{k=0}^{K-1} c_k \Psi_k(\mathbf{x}) = \boldsymbol{\Psi}(\mathbf{x})^T \mathbf{c} \in \mathcal{GP}(\boldsymbol{\Psi}(\mathbf{x})^T \boldsymbol{\mu}, \boldsymbol{\Psi}(\mathbf{x}) \boldsymbol{\Sigma} \boldsymbol{\Psi}(\mathbf{x}')^T)$$

# Sensitivity information comes free with PC surrogate,

$$g(x_1, \dots, x_d) = \sum_{k=0}^{K-1} c_k \Psi_k(\mathbf{x})$$

- Main effect sensitivity indices

$$S_i = \frac{Var[\mathbb{E}(g(\mathbf{x}|x_i)]}{Var[g(\mathbf{x})]} = \frac{\sum_{k \in \mathbb{I}_i} c_k^2 \|\Psi_k\|^2}{\sum_{k>0} c_k^2 \|\Psi_k\|^2}$$

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- Joint sensitivity indices

$$S_{ij} = \frac{Var[\mathbb{E}(g(\mathbf{x}|x_i, x_j)]}{Var[g(\mathbf{x})]} - S_i - S_j = \frac{\sum_{k \in \mathbb{I}_{ij}} c_k^2 \|\Psi_k\|^2}{\sum_{k>0} c_k^2 \|\Psi_k\|^2}$$

$\mathbb{I}_{ij}$  is the set of bases with only  $x_i$  and  $x_j$  involved

Sensitivity information comes free with PC surrogate,  
but not with piecewise PC

$$g(x_1, \dots, x_d) = \sum_{k=0}^{K-1} c_k \Psi_k(\mathbf{x})$$

- Main effect sensitivity indices

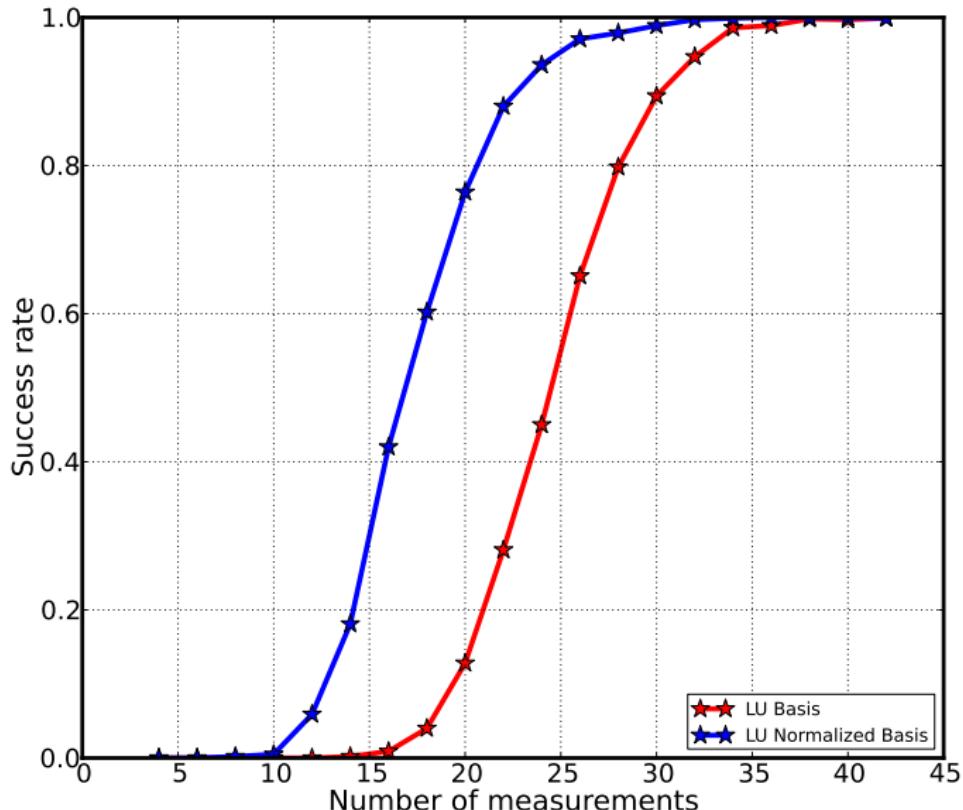
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- For piecewise PC, need to resort to Monte-Carlo estimation  
[Saltelli, 2002].

# Basis normalization helps the success rate



## Input correlations: Rosenblatt transformation

- Rosenblatt transformation maps any (not necessarily independent) set of random variables  $\lambda = (\lambda_1, \dots, \lambda_d)$  to uniform i.i.d.'s  $\{x_i\}_{i=1}^d$  [Rosenblatt, 1952].

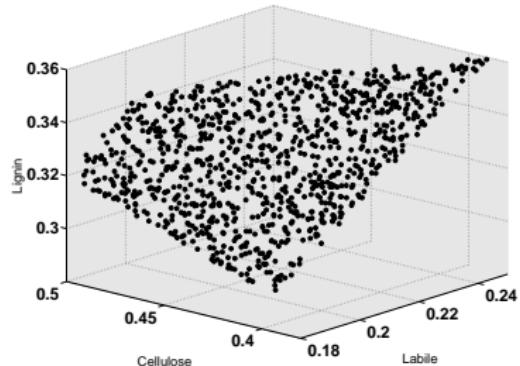
$$x_1 = F_1(\lambda_1)$$

$$x_2 = F_{2|1}(\lambda_2|\lambda_1)$$

$$x_3 = F_{3|2,1}(\lambda_3|\lambda_2, \lambda_1)$$

$$\vdots$$

$$x_d = F_{d|d-1, \dots, 1}(\lambda_d|\lambda_{d-1}, \dots, \lambda_1)$$



- Inverse Rosenblatt transformation  $\lambda = R^{-1}(\mathbf{x})$  ensures a well-defined input PC construction

$$\lambda_i = \sum_{k=0}^{K-1} \lambda_{ik} \Psi_k(\mathbf{x})$$

- Caveat: the conditional distributions are often hard to evaluate accurately.

# Strong discontinuities/nonlinearities challenge global polynomial expansions

- Basis enrichment [Ghosh & Ghanem, 2005]
- Stochastic domain decomposition
  - Wiener-Haar expansions,  
Multiblock expansions,  
Multiwavelets, [Le Maître *et al*, 2004,2007]
  - also known as Multielement PC [Wan & Karniadakis, 2009]
- Smart splitting, discontinuity detection  
[Archibald *et al*, 2009; Chantrasmi, 2011; Sargsyan *et al*, 2011; Jakeman *et al*, 2012]
- Data domain decomposition,
  - Mixture PC expansions [Sargsyan *et al*, 2010]
- Data clustering, classification,
  - Piecewise PC expansions

## Piecewise PC expansion with classification

- Cluster the training dataset into non-overlapping subsets  $\mathcal{D}_1$  and  $\mathcal{D}_2$ , where the behavior of function is smoother
- Construct global PC expansions  $g_i(\mathbf{x}) = \sum_k c_{ik} \Psi_k(\mathbf{x})$  using each dataset individually ( $i = 1, 2$ )
- Declare a surrogate

$$g_s(\mathbf{x}) = \begin{cases} g_1(\mathbf{x}) & \text{if } \mathbf{x} \in^* \mathcal{D}_1 \\ g_2(\mathbf{x}) & \text{if } \mathbf{x} \in^* \mathcal{D}_2 \end{cases}$$

\* Requires a classification step to find out which cluster  $\mathbf{x}$  belongs to. We applied Random Decision Forests (RDF).

- Caveat: the sensitivity information is harder to obtain.