Uncertainty quantification in high-dimensional, nonlinear and expensive models



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- The case for Uncertainty Quantification (UQ): what, why and how
- Surrogates needed for expensive models
- Polynomial Chaos (PC) spectral methods for forward UQ
- Bayesian compressive sensing finds sparse surrogates in high-d
- Classification ideas deal with nonlinearities

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 - Classification and piecewise-PC deal with nonlinearities

Laundry list

Uncertainty sources

- Model parameters
- Initial/boundary conditions
- Model geometry/structure
- Lack of knowledge
- Data noise
- Intrinsic stochasticity
- Numerical errors
- Faults, data loss

UQ enables

- Model predictions with uncertainty
- Model validation and comparison
- Confidence assessment
- Dimensionality reduction
- Optimal design
- Reliability analysis
- Decision support
- Data assimilation

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- (Bayesian) Probabilistic UQ methods well-suited
- Uncertainty \sim Probability \sim State of knowledge









Application of Interest: Community Land Model



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

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

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UQ challenges:

- Computationally expensive
- High dimensionality
- Non-smooth/nonlinear behavior

Surrogates are necessary for computationally expensive models

Construct surrogate for a complex model f(x) to enable sampling-intensive studies:

- Global sensitivity analysis
- Optimization
- Uncertainty propagation (Forward UQ)
- Input parameter calibration (Inverse UQ)
- • •
- 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
 - Too many parameters in the surrogate parameterization

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Polynomial Chaos surrogate construction

• Scale the input parameters $x_i \in [a_i, b_i]$

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

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

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

- A lot of information for free:
 - Global sensitivity (Sobol) indices, variance-based decomposition
 - Moments of *u*, as a random variable

Alternative methods to obtain PC coefficients

$$y = f(x) \simeq \sum_{k=0}^{K-1} c_k \Psi_k(\boldsymbol{x})$$

- <u>Projection</u> $c_k = \frac{\langle f(\boldsymbol{x})\Psi_k(\boldsymbol{x})\rangle}{\langle \Psi_k^2(\boldsymbol{x})\rangle}$ The integral $\langle f(\boldsymbol{x})\Psi_k(\boldsymbol{x})\rangle = \int f(\boldsymbol{x})\Psi_k(\boldsymbol{x})d\boldsymbol{x}$ can be estimated by
 - Monte-Carlo



Quadrature



many(!) random samples



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samples at quadrature

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Bayesian regression

 $P(c_k|f(\mathbf{x}_j)) \propto P(f(\mathbf{x}_j)|c_k)P(c_k)$







any (number of) samples

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Posterior



many(!) random samples



samples at quadrature

any (number of) samples

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

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

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

$$\mathbf{y} \approx \mathbf{Pc}$$
 or $y_n \approx \sum_k c_k \Psi_k(\mathbf{x}_n)$

- 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 ⇒ need some regularization/constraints.
- Least-squares $\operatorname{argmin}_{\boldsymbol{c}} \{ || \boldsymbol{y} \boldsymbol{P} \boldsymbol{c} ||_2 \}$
- The 'sparsest' $argmin_{c} \{ ||y Pc||_{2} + \alpha ||c||_{0} \}$
- Compressive sensing $\operatorname{argmin}_{\boldsymbol{c}} \{ || \boldsymbol{y} \boldsymbol{P} \boldsymbol{c} ||_2 + \alpha || \boldsymbol{c} ||_1 \}$

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 Bayesian

```
\begin{array}{c} \textit{argmin}_{\pmb{c}} \left\{ ||\pmb{y} - \pmb{P} \pmb{c}||_2 + \alpha ||\pmb{c}||_1 \right\} \\ \text{Likelihood} \quad \text{Prior} \end{array}
```

BCS removes unnecessary basis terms



The square (i, j) represents the (log) spectral coefficient for the basis term $\psi_i(x_1)\psi_j(x_2)$.

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Success rate grows with more data and 'sparser' model

Consider test function

$$f(\boldsymbol{x}) = \sum_{k=0}^{K-1} c_k \Psi_k(\boldsymbol{x})$$

where only S coefficients c_k are non-zero. Typical setting is

S < N < K



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Recovering true PC coefficients



Recovering true PC coefficients



Recovering true PC coefficients



Basis set growth with iterative BCS





Piecewise-PC expansion deals with nonlinearities:

use data classification methods

- Cluster the training dataset into non-overlapping subsets D₁ and D₂, 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

$$g_s(oldsymbol{x}) = egin{cases} g_1(oldsymbol{x}) & ext{if } oldsymbol{x} \in^* \mathcal{D}_1 \ g_2(oldsymbol{x}) & ext{if } oldsymbol{x} \in^* \mathcal{D}_2 \end{cases}$$

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

• Caveat: the sensitivity information is harder to obtain.





Piecewise 2-nd order surrogate



Piecewise 5-th order surrogate



Piecewise 5-th order surrogate



Piecewise 5-th order surrogate





Sparse PC surrogate for the Community Land Model

- Main effect sensitivities : rank input parameters
- Joint sensitivities : most influential input couplings
- About 200 polynomial basis terms in the 70-dimensional space
- Sparse PC will further be used for
 - · sampling in a reduced space
 - parameter calibration against experimental data





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Summary

- Surrogate models are necessary for UQ studies of 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 *iterative* Bayesian CS algorithm that reduces dimensionality and increases order on-the-fly.
- Nonlinear behavior
 - Data clustering and classification-driven piecewise PC
- Future work, open issues
 - Computational design. What is the best sampling strategy?
 - Weighted *l*₁ minimization to accomodate natural coefficient decay.

Literature

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Thank You

Random variables represented by Polynomial Chaos

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

• $\boldsymbol{\xi} = (\xi_1, \cdots, \xi_d)$ standard i.i.d. r.v. Ψ_k standard polynomials, orthogonal w.r.t. $\pi(\boldsymbol{\xi})$.

 $\Psi_k(\xi_1,\xi_2,\ldots,\xi_d) = \psi_{k_1}(\xi_1)\psi_{k_2}(\xi_2)\cdots\psi_{k_d}(\xi_d)$

- Typical truncation rule: total-order p, $k_1 + k_2 + ... k_d \le p$. 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, <u>Legendre-Uniform</u>, (discrete) Poisson-Charlier.

Bayesian Compressive Sensing (BCS), or Relevance Vector Machine (RVM)

Dimensionality reduction by using hierarchical priors

$$p(c_k|\sigma_k^2) = \frac{1}{\sqrt{2\pi}\sigma_k} e^{-\frac{c_k^2}{2\sigma_k^2}} \qquad \qquad p(\sigma_k^2|\alpha) = \frac{\alpha}{2} e^{-\frac{\alpha\sigma_k^2}{2}}$$

Effectively, one obtains Laplace sparsity prior

$$p(\boldsymbol{c}|\alpha) = \int \prod_{k=0}^{K-1} p(c_k|\sigma_k^2) p(\sigma_k^2|\alpha) d\sigma_k^2 = \prod_{k=0}^{K-1} \frac{\sqrt{\alpha}}{2} e^{-\sqrt{\alpha}|c_k|}$$

- The parameter α can be further modeled hierarchically, or fixed.
- Evidence maximization dictates values for σ²_k, α, σ² and allows exact Bayesian solution

$$m{c} \sim \mathcal{MVN}(m{\mu}, m{\Sigma})$$

with

$$\boldsymbol{\mu} = \sigma^{-2} \boldsymbol{\Sigma} \boldsymbol{P}^{T} \boldsymbol{u} \qquad \boldsymbol{\Sigma} = \sigma^{2} (\boldsymbol{P}^{T} \boldsymbol{P} + \text{diag}(\sigma^{2}/\sigma_{k}^{2}))^{-1}$$

• KEY: Some $\sigma_k^2 \rightarrow 0$, hence the corresponding basis terms are dropped.

[Tipping, 2001, Ji et al., 2008; Babacan et al., 2010]

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.* 2013].



Input correlations: Rosenblatt transformation

Rosenblatt transformation maps any (not necessarily independent) set of random variables λ = (λ₁,..., λ_d) to uniform i.i.d.'s {η_i}^d_{i=1}
 [Rosenblatt, 1952].





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

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

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

Sensitivity information comes free with PC surrogate,

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

Main effect sensitivity indices

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

 I_i is the set of bases with only x_i involved

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Joint sensitivity indices

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

 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

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

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