## Uncertainty quantification in computer experiments with polynomial chaos

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Uncertainty quantification (UQ) in computer experiments

- $\triangleright$  Context: Deterministic and complex numerical simulator are used to model real dynamic systems and they can be computationally expensive to run
- $\triangleright$  We are interested to study the effect of epistemic (lack of knowledge) and aleatoric (inherent to system) uncertainties on the model outputs
- $\triangleright$  Sources include initial condition, boundary condition  $\&$  model parameters
- $\triangleright$  Example: drug clearance in circulation as an exponential decay response  $\frac{d\theta}{dt} = -C\theta$  with  $C$  as a r.v. that represents the population response
- $\triangleright$  Conventional approaches such as MC are not practical in studying these expensive simulators
- Goal: PC construct a metamodel that mimics the complex model's behaviour and conduct UQ, SA, quantile estimation, optimization, calibration, etc.

## Probabilistic framework

The UQ of a computer experiment follows the following iterative steps:

- 1. representation of input uncertainties random variable or process
- 2. uncertainty propagation MC, GP or gPC
- 3. quantification of solution uncertainty mean, variance, pdf or sensitivity



De Rocquigny (2006)

#### Stochastic input representation: stochastic process

Any second order random process  $\kappa(x, \omega)$ , with continuous and bounded covariance kernel  $C(x_1, x_2) = \mathbb{E}(\kappa(x_1, \omega) \otimes \kappa(x_2, \omega))$ , can be represented as an infinite sum of random variables. It is real, symmetric and positive–definite.



 $\triangleright$  Karhunen-Loève (KL) expansion represents the random process with an orthogonal set of deterministic functions with random coefficients as

$$
\kappa(x,\omega)=\mu_{\kappa}(x)+\sum_{n=1}^N\sqrt{\lambda_n}\psi_n(x)\xi_n(\omega).
$$

- $\triangleright$  For a continuous kernel, the convergence of the KL expansion is uniform as  $N \rightarrow \infty$ . Karhunen (1948) & Loève (1977)
- $\blacktriangleright \psi_n(x)$  and  $\lambda_n$  solved from Fredholm integral equation of 2nd kind with  $C(x_1, x_2)$ .

Stochastic input representation: random variables

**F** Represent the random variable,  $\kappa(\omega)$ , with orthogonal functions of the stochastic variable with deterministic coefficients

$$
\kappa(\omega)=\sum_{m=0}^{\infty}\kappa_m\phi_m(\xi(\omega)).
$$

- $\triangleright$  Wiener-Chaos: representation of a Gaussian random variable using Hermite polynomials with  $L^2$  convergence as  $M\to\infty$ . Wiener (1938), Ghanem & Spanos (1991) and Cameron & Martin (1947)
- $\triangleright$  generalized Polynomial Chaos: generalized representation to non-Gaussian random variables with polynomials from the Wiener–Askey scheme. Xiu & Karniadakis (2002)
- If  $\kappa(\omega)$  follows a normal distribution, it can be represented exactly as  $\kappa(\omega) = \mu_{\kappa} + \sigma_{\kappa} \xi$  where  $\xi$  is the linear term in Hermite

## Selection of orthogonal basis

In the propagation step, we need to evaluate the inner product w.r.t. the probability space measure,  $\rho(\xi)d\xi$  as

$$
\langle \phi_i(\xi), \phi_j(\xi) \rangle = \int_{\Gamma} \phi_i(\xi) \phi_j(\xi) \rho(\xi) d\xi.
$$

► Correspondence between the *pdf* of  $\xi$ ,  $\rho(\xi)$ , and the weighting function of classical orthogonal polynomials,  $w(\xi)$ , determines the polynomial basis



## Multivariate basis

Multivariate basis is the tensor products of 1D polynomials

$$
\phi_m(\xi) = \phi^{\alpha_{m,n=1}}(\xi_1) \otimes \phi^{\alpha_{m,n=2}}(\xi_2) \otimes \cdots \otimes \phi^{\alpha_{m,n=N}}(\xi_N), \text{ for } m = 0, \cdots, M,
$$
  
=  $\phi^{\alpha_m}(\xi)$ , for  $m = 0, \cdots, M$ .

Truncation depends on input dimension, N, and output nonlinearity, P





## Stochastic Galerkin method: intrusive approach

PC represent the stochastic solution  $u(x, \xi)$  with the same orthogonal basis as the input, *i.e.*  $u(\mathbf{x}, \xi) = \sum u_m(\mathbf{x}) \phi_m(\xi)$ Substitute the expansions into the system of equations,  $\mathcal{L}(\mathbf{x}, \boldsymbol{\xi}; u) = f(\mathbf{x}, \boldsymbol{\xi})$ . Take the Galerkin projection, *i.e.* 

$$
\langle \mathcal{L}(\mathbf{x}, \xi; \sum u_m(\mathbf{x}) \phi_m(\xi)) \rangle, \phi_m(\xi) \rangle = \langle f(\mathbf{x}, \xi), \phi_m(\xi) \rangle, \text{ for } m = 0, ..., M.
$$

- $u_m(x)$  are solved from the system of  $(M + 1)$  coupled equations.
- $\triangleright$  The system is deterministic and can be solved using a standard discretization technique.
- $\triangleright$  Extensive modification on the simulator is needed.

## Stochastic Galerkin method: intrusive approach

#### Example

First-order linear ODE:  $\Theta(t,\xi) = -C(\xi)\Theta(t,\xi)$  with rate of decay as a normal r.v., *i.e.*  $C(\xi)=\sum_{i=0}^{M}C_{i}\phi_{i}(\xi).$  The gPC expansions of  $C(\xi)$  and  $\Theta(t,\xi)$  are substituted into the ODE to give

$$
\sum_{k=0}^{M_{\theta}}\dot{\Theta}_k(t)\phi_k(\xi) = -\sum_{i=0}^{M_C}\sum_{j=0}^{M_{\theta}}C_i\Theta_j(t)\phi_i(\xi)\phi_j(\xi).
$$

The Galerkin projection of the expanded ODE with orthogonal polynomial:

$$
\dot{\Theta}_k(t)=-\sum_{i=0}^{M_C}\sum_{j=0}^{M_\theta}\frac{\langle\phi_i\phi_j\phi_k\rangle}{\langle\phi_k^2\rangle}C_i\Theta_j(t), \text{ for } k=0, ..., M_\theta.
$$

This coupled deterministic system of equations is solved with an initial condition  $\Theta(t=0) = \sum \Theta_m(t=0) \phi_m(\xi)$ . With increasing t, the modal coefficients are propagated from the lower  $\Theta_m$  to higher  $\Theta_m$ , *i.e.* propagation of uncertainty as increasing non–linear response in the random space.

## Surface response of the linear ODE

- $\dot{\Theta}(t,\xi) = -C(\xi)\Theta(t,\xi)$
- $\Theta(t, \xi)$  response is exponential in t with  $\Theta(t = 0) = 1$ .
- **Figure 1** Treating the coefficient of decay as a random variable,  $C(\xi) \sim \mathcal{N}(1, 1)$
- $\triangleright$  We represent the univariate stochastic output  $\Theta(t;\xi)$  as a linear combination of Hermite polynomials  $\Theta(t;\xi) = \sum \Theta_m(t) \phi_m(\xi)$ .
- $\triangleright$  Uncertainty propagation visualized as solution response surface evolution in random space,  $\xi$



#### The choice of polynomial chaos truncation

- As response in  $\xi$  becomes more non-linear with t, the higher order P in  $\phi_m(\xi)$  are needed in gPC expansion
- $\triangleright$  Estimation of higher order statistics also require higher P
- $\triangleright$  Premature truncation leads to large error in the response surface and the solution statistics



## Evolution of the PC coefficients

Increasing t propagates the initial uncertainty from lower order coefficients to higher order coefficients



- **I** The task now is to determine the coefficients of expansion,  $\Theta_m(t)$  in the representation.
- $\triangleright$  This simple system of equation easily solved with the intrusive approach
- $\triangleright$  Complex numerical solvers can benefit from a non–intrusive approach

Projecting directly the stochastic solution,  $u(\mathbf{x}, \boldsymbol{\xi}) = \sum u_m(\mathbf{x}) \phi_m(\boldsymbol{\xi})$ , onto the orthogonal basis,  $\phi_m(\xi)$ , we obtain the following  $(M + 1)$  decoupled equations:

$$
u_m(\mathbf{x}) = \frac{\langle u(\mathbf{x}, \boldsymbol{\xi}), \phi_m(\boldsymbol{\xi}) \rangle}{\langle \phi_m^2(\boldsymbol{\xi}) \rangle}, \quad \text{for } m = 0, ..., M.
$$

The inner–product can be evaluated using Monte Carlo or related methods. We investigate a numerical quadrature approach to approximate the inner product where the numerical solver is treated as a black box from which samples are repeated taken.

#### One–dimensional quadrature rules

Integrals are approximated as the weighted sum of function evaluations on deterministic quadrature points, *i.e.* 

$$
\langle u(\mathbf{x},\boldsymbol{\xi}),\phi_m(\boldsymbol{\xi})\rangle = \int_{\Gamma} u(\mathbf{x},\boldsymbol{\xi})\phi_m(\boldsymbol{\xi})\rho(\boldsymbol{\xi})d\boldsymbol{\xi},
$$

$$
\approx \sum_{j=0}^{N_q} w_j u(\mathbf{x},\mathbf{z}_j)\phi_m(\mathbf{z}_j).
$$

The accuracy of the method depends on the selection of the quadrature approach, *i.e.* constructions of  $w_i$  and  $z_j$ .



Multi–dimensional quadrature rules are constructed from 1D quadrature rules.

#### Full–tensor quadrature

Multi–dimensional full–tensor quadrature relies on tensor product of 1D quadrature rules, e.g. N–dimensional quadrature points are

$$
\mathcal{Q}_L^N(f)=(\mathcal{U}^{i_1}\otimes\cdots\otimes\mathcal{U}^{i_N})(f).
$$

Example: Two–dimensional Gauss–Legendre quadrature:



**Accuracy:** Theoretical polynomial exactness  $P = 2L - 1$  in each dimension where L is the number of quadrature points in each dimension **Cost:** Number of quadrature points grows as  $\mathcal{O}(L^N)$  and error converges as  $\epsilon(Z) = \mathcal{O}(Z^{-r/N})$  – "curse of dimensionality"

## Sparse quadrature: the Smolyak approach

"Curse of dimensionality" could be 'broken' with the sparse grid. Its construction is based on the following three steps: Gerstner & Griebel (1998)

- 1. Constructed from 1D difference grid
- 2. Tensor product of 1D difference grids: cost reduction
- 3. Linear combination of the tensor products: embeddedness  $\rightarrow$  refinement cost reduction

**Accuracy:** Theoretical polynomial exactness at least  $P \leq 2L - 1$  where L is the quadrature level. Smolyak (1963), Novak & Ritter (1996) **Cost:** Error converges as  $\epsilon(Z) = \mathcal{O}(Z^{-r}(\log(Z)^{(N-1)(r+1)}))$ . Novak & Ritter (1996)

Sparse quadrature: with nested Clenshaw-Curtis quadrature rule

 ${\bf 1D}$  difference grid:  $\triangle^1_k f := \left( {\cal Q}^1_k - {\cal Q}^1_{k-1} \right) f$ 



Sparse quadrature: with nested Clenshaw-Curtis quadrature rule

**1D difference grid:** 
$$
\triangle_k^1 f := (Q_k^1 - Q_{k-1}^1) f
$$
 **Tensor product:**  $(\triangle_{k_1}^1 \otimes \cdots \otimes \triangle_{k_N}^1) f$ 



Sparse quadrature: with nested Clenshaw-Curtis quadrature rule

 ${\bf 1D}$  difference grid:  $\triangle^1_k f := \left( {\cal Q}^1_k - {\cal Q}^1_{k-1} \right) t$ Tensor product:  $\left(\triangle^1_{k_1} \otimes \cdots \otimes \triangle^1_{k_N}\right)$  / Linear combination:  $\mathcal{Q}_L^N[f] := \sum \left(\bigtriangleup^1_{k_1}\otimes\cdots\otimes\bigtriangleup^1_{k_N}\right)$   $t$ 



#### Sparse quadrature: comparison with full–tensor quadratures



Sparse Clenshaw-Curtis Chebyshev: P=7, P=9 & P=11

Full Gauss-Legendre Quadrature: P=7, P=9 & P=11



#### Canonical, maximum and anisotropic expansions

 $M$  is determined by the accuracy of the quadrature approach. If the quadrature has a polynomial accuracy of  $P$  or  $P$ , there are the following expansions for

$$
f_r(\mathbf{x}) = \sum_{\alpha \in \mathbb{N}^N} f_\alpha \phi_\alpha(\mathbf{x})
$$

- **►** Canonical: total degrees not greater than P, i.e.  $\{\phi_{\alpha} / |\alpha| \leq P\}$
- **Maximum:** degree in each n not greater than P, i.e.  $\{\phi_{\alpha} \mid \alpha \leq P\}$ .
- Anisotropic: degree in each n not greater than  $P_n$ , i.e.  $\{\phi_{\alpha} \mid \alpha \leq P\}$ .



gPC as a Uncertainty Quantification (UQ) & Sensitivity Analysis (SA) tool

Statistical moments:

$$
\mu_u(\mathbf{x}) = \int_{\Gamma} u_r(\mathbf{x}; \omega) \phi_0(\xi) \rho(\xi) d\xi = u_0(\mathbf{x}),
$$
  

$$
\sigma_{u, gPC}^2(\mathbf{x}) = \int_{\Gamma} \left[ \sum_{m=0}^M u_m(\mathbf{x}) \phi_m(\xi) - u_0(\mathbf{x}) \right]^2 \rho(\xi) d\xi = \sum_{m=1}^M u_m^2(\mathbf{x}) \langle \phi_m^2(\xi) \rangle.
$$

**Solution sensitivity:** Partial differentiation wrt  $\xi_n$  Agarwal (2008)

$$
\mathbf{S}_{\xi_n}(\mathbf{x}) = \frac{\partial u_r(\mathbf{x}; \boldsymbol{\xi})}{\partial \xi_n}
$$

.

Sensitivity analysis: partial variances Sobol' (1993)

$$
\sigma^2_u(\mathbf{x}) = \sum_{i_1=1}^N D_{i_1}(\mathbf{x}) + \sum_{i_1=1}^N \sum_{i_2=1}^{i_1} D_{i_1 i_2}(\mathbf{x}) + \sum_{i_1=1}^N \sum_{i_2=1}^{i_1} \sum_{i_3=1}^{i_2} D_{i_1 i_2 i_3}(\mathbf{x}) \cdots + D_{i_1 i_2 \ldots i_N}(\mathbf{x}).
$$

Probability density function (PDF): numerical computation from the histogram of a large MC sample of  $u_r(x,\xi)$  based on the distribution of  $\xi$ 

# Application of gPC to some examples



## Sensitivity of spatially developing mixing layer

- $\triangleright$  Coherent vortical structures triggered by inflow forcing Brown & Roshko (1974)
- **In Shear layer at the inflow approximated as**  $\overline{U}_{in}(y) = 1 + \lambda \tanh(y/2)$
- $\triangleright$  Downstream shear layer growth is very sensitive to forcing definition
- $\triangleright$  Forcing with LST fundamental mode, *i.e.* most unstable, and its subharmonic modes:  $u_p(y, t) = \sum \epsilon_n f_n(y) \exp(i(\omega_n t + \gamma_n))$
- ▶ 3D flow structure is largely  $2D \rightarrow 2D$  DNS Delville et al. (1999)
- $\blacktriangleright$  Goal: To generalize the approach to design discrete forcing with random magnitude or phasing



## Sensitivity to forcing: magnitude  $\epsilon_n$

- $\blacktriangleright$  Instantaneous vorticity contours with bimodal perturbation
- $\triangleright$  Vortical structure variation as the relative frequency content in inflow forcing changes



## Stochastic mixing layer with random magnitudes,  $\epsilon_n$

Treat  $\epsilon_n$  and  $\gamma_n$  as random variables to determine the most general way to control mixing layer growth with inflow forcing.

- $\triangleright$  Bimodal forcing and trimodal forcing examined
- Stochastic forcing magnitudes  $\epsilon_n$  as uniform variables in [0, 10% $\overline{U}$ ]
- $\blacktriangleright$  Legendre-Chaos expansion of stochastic fields
- $\triangleright$  Mixing layer solutions with 2D spectral/hp DNS solver
- $\blacktriangleright$  Re = 100,  $\lambda = 0.5$
- $\triangleright$  Non–intrusive Probabilistic Collocation Method with full–tensor Gauss–quadrature
- ▶ 81 full–tensor quadrature points for bimodal forcing (N=2, L=9, P=8) & 1000 for trimodal  $(N=3, L=10, P=9)$
- Examine time-averaged mixing layer thickness, e.g. momentum thickness  $\theta$

## Accuracy of the gPC expansion: solution prediction

With  $u(x,\xi) = u_m(x)\phi_m(\xi)$ , we can predict the solution at an arbitrary point within Γ. Accuracy of the prediction increases with increasing *M* or *P*.



#### Response variability in trimodal perturbation case

- Initial response up to  $x/\theta_{in} = 250$  similar to the bimodal case
- $\blacktriangleright$  Large local variance at the location associated with the onset of deterministic subharmonic vortex merging



## Partial variance contour in trimodal vorticity



- $\blacktriangleright$  D<sub>n</sub>: sensitivities of the solution to  $\epsilon_n$
- $\blacktriangleright$  Contours of each sensitivity index correspond closely to the deterministic vortex-roll up of each mode

#### Partial variance in trimodal vorticity contour



- interaction between  $\epsilon_i$  and  $\epsilon_i$
- $10^{100}$ <br>  $10^{100}$ <br> ► Large  $D_{12}$  and  $D_{23} \rightarrow$  interactions between successive modes are dominant Kelly (1967)
	- $\triangleright$   $D_{123}$ : sensitivities of the solution to the mutual interaction amongst all modes

## $\theta$  PDF in trimodal perturbation case



## Stochastic mixing layer with random phase  $\gamma_n$

- $\triangleright$  Bimodal forcing and trimodal forcing examined
- Stochastic phase shifts  $\gamma_n$  as uniform random variables in [0, 2π)
- Forcing magnitudes maintained at  $\sum \epsilon_n = 10\% \overline{U}$
- $\triangleright$  SCM with Newton-Cotes quadrature
- $\triangleright$  Fourier-Chaos expansion of stochastic fields
- $\triangleright$  Discrete Fourier transformation (DFT) speeds up coefficient computations
- $\triangleright$  72 equidistant quadrature samples are used (nested points)
- Examine time-averaged mixing layer thickness, e.g. momentum thickness  $\theta$

### Response of momentum thickness



- Symmetry observed as  $\gamma_2 \in [0, 2\pi]$ includes two periods of fund. forcing
- $\triangleright$  Mixing layer growth strongly delayed over small  $\gamma_2$  range near  $70^\circ$  Inoue (1995)
- Delayed growth reported for  $\gamma_2 = 0$  at merging locations Stanley & Sarkar (1997)
- ▶ 45° difference between inflow forcing formulations
- Phase shift at inflow does not correspond to phase shift at merging locations

#### Mixing layer growth rate statistics





- $\triangleright$  'Normal growth': Fast growth near inflow followed by sharp drop in  $\partial\theta/\partial x$ . Drop or contraction of the mixing layer Oster & Wygnansk (1982)
- $\triangleright$  'Delayed growth': Slower growth with less  $\partial\theta/\partial x$  fluctuation. Large variance due to solution sensitivity in  $\gamma_2 \in [45^\circ, 80^\circ]$ . Range of sensitivity is small Stanley & Sarkar (1997)

## PC as a quantile estimation tool

**Empirical quantile**: estimated from  $\hat{Y}_\alpha = \inf\{y : \hat{F}(y) > \alpha\}$  which gives

$$
\hat{Y}_{\alpha} = Y_{(\lceil \alpha Z \rceil)},\tag{1}
$$

where  $\{Y_{(i)}\}_{i=1}^Z$  are the ordered set of the Z MC samples. The metamodel accurately determines the statistical moments but fails in extreme quantile estimations, i.e.  $\alpha$  near 0 or 1. We propose a multi-element refinement approach: global gFC metamodel is complimented by local metamodel constructed around design points  $\xi_{\alpha}$ . **Design point**: *most likely* random input that corresponds to  $u_r(\mathbf{x}, \xi) = u_\alpha(\mathbf{x})$ . This gives a constraint nonlinear minimization problem, *i.e.* 

$$
\min ||\xi||, \quad \text{s.t.} \quad \sum_{m=0}^{M} u_m(\mathbf{x}) \phi_m(\xi) - \hat{Y}_{\alpha} = 0.
$$

The above problem is solved by the method of **Lagrangian multipliers**.

#### Multi–Element Monte Carlo simulation

Local gPC metamodels are created around the design points. The multielement solution is used as the metamodel, *i.e.* 

$$
D_{\text{ME}} = \begin{cases} D_{\text{global}} = D \setminus D_{\text{local}}, & \text{domain of global gPC}, \\ D_{\text{local}} = \cup D_{\beta_i}, & \text{domains of refinement about } \hat{\xi}_{\alpha_i}, \text{ for } i = 1, ..., N_{\beta}. \end{cases}
$$



The final multi–element gPC (MEgPC) metamodel is

$$
f_{\text{ME}}(\mathbf{x}) = \begin{cases} \sum_{m=0}^{M} f_m \phi_m(\mathbf{x}), & \text{if } \mathbf{x} \in D_{\text{global}},\\ \sum_{m=0}^{M_i^*} f_{m,i}^* \psi_{m,i}(\mathbf{T}_i^{-1}(\mathbf{x})), & \text{if } \mathbf{x} \in D_{\beta_i}. \end{cases}
$$

where  $\mathbf{T}_i$  a transformation operator that maps a point in the uniform bounded support  $\mathbf{x}^* \in [-1,1]^N$  to the local domain  $\mathbf{x} \in D_{\beta_i}.$ 

#### Example: Gaussian–like response

We examine the quantile of the output of a Gaussian-like function:

$$
f(\mathbf{x}) = \sum_{i=1}^{N_{\alpha}} \prod_{n=1}^{N} \exp\left(\frac{-\left(x_n - \mu_{n,i}\right)^2}{2\sigma_{n,i}^2}\right),\tag{2}
$$

where  $\|\mu\| = 2$ ,  $\sigma = 1$ , x are i.i.d. random variables and  $x_n \in \mathcal{N}(0, 1)$ .

# **Multi–element Metamodel**<br>Quadratic global metamodel with quadratic local metamodel



 $\alpha$ –quantile estimator convergence for MC, IS and global gPC

 $\blacktriangleright$  Monte Carlo  $\hat{Y}_\alpha$  converges as  $1/\sqrt{Z}$ 

- Importance sampling  $\hat{Y}_{\alpha}$  computed at selected Z: Z/2 MC samples for first estimate of  $\hat{Y}_{\alpha}$ , at most  $Z/4$  for GPM and the rest for IS
- ► Global full and sparse gPC estimations of  $\hat{Y}_{\alpha,r}$  (from  $L = 3$  to 7) are poor



## Effects of different local refinements

- ► Local full (canonical & maximum) and sparse gPC metamodel refinements
- **I** Maximum expansion improves the accuracy of  $\hat{Y}_{\alpha,ME}$  given the same Z
- $\blacktriangleright$  Seek best  $\hat{\xi}_{\alpha,\tau}$  estimation by maximizing  $Z$  in global gPC metamodel



#### Target cost study

- An arbitrary target cost that increases linearly with N:  $Z_{total} = 100N$
- $\triangleright$  Monte Carlo and importance sampling  $\hat{Y}_{\alpha}$  with entire sampling budget
- Global full and sparse + local full maximum  $(+)$  and sparse  $(°)$ supplemental metamodels
- $\triangleright$  Maximize global metamodel cost while not exceeding the entire budget



## Example: Hypertangent response

We examine the quantile of the output of a hypertangent function:

$$
Y(\mathbf{x}) = 1 + \tanh\left(\sum_{n=1}^N \sigma_n(x_n - \mu_n)\right).
$$

where the N–dimensional input are i.i.d. random variables  $x \in \mathcal{N}(0, 1)$ .



## Anisotropic grid

- $\blacktriangleright$  The dominance of some random variables can be revealed by examining the partial variance of the global gPC metamodel
- $\blacktriangleright$  One–dimensional metamodels about  $\hat{\xi}_{\alpha,r}$  can identify dominant directions



Anisotropic grids, P in  $\hat{\xi}'_c$  $\alpha_{\alpha,r}^{\top}$  and linear in transverse directions, reduce cost



#### Target cost study

- An arbitrary target cost that increases linearly with  $N: Z_{total} = 100N$
- $\triangleright$  Monte Carlo and importance sampling  $\hat{Y}_{\alpha}$  with entire sampling budget
- Global full and sparse + local full canonical  $(\Box)$  and anisotropic  $(\triangle)$ supplemental metamodels
- $\blacktriangleright$  Maximize global metamodel cost while not exceeding the entire budget



#### Quantile of multivariate output

We assume that all components of the random output Y are extreme and define the multivariate  $\alpha$ –quantile as the point y<sub>α</sub> where the multivariate and marginal cdf's satisfy the following conditions

$$
F(\mathbf{y}_{\alpha}) = \alpha \quad \text{and} \quad F_1(y_{\alpha,1}) = F_2(y_{\alpha,2}) = \cdots = F_K(y_{\alpha,K}) \tag{3}
$$

where K is the number of outputs. Results with N=2 and  $\alpha = 99\%$  case for multiple Gaussian peaks:



## Calibration and sensitivity analysis GCM

- $\triangleright$  Examine the AGCM ECHAM6 with uncertain parameters in cloud modeling
- $\rightarrow$  1977 climatological distributions of sea ice and surface temperature used as initial condition
- $\blacktriangleright$  Five R.V. in the expert range transformed to the Gaussian space
- $\triangleright$  Ensemble of model output created for a single year run
- $\blacktriangleright$  Full–tensor quadrature with squadratic accuracy, *i.e.* 243 points

## Selection of the input random variables

Table: Expert parameter range and their default values



- $\triangleright$  zinhomi & zinhoml are treated as uniform r.v.
- $\triangleright$  entrscv, entrpen & cprcon are treated as uniform r.v or log uniform r.v.
- A dependent parameter,  $\textit{cmfctop} = \textit{entrscv} \times \frac{1000}{3}$ , is included
- $\triangleright$  A uniform distribution under–weights the entire lower range



#### Validation: Comparison of computed global contours and gPC predictions

- $\triangleright$  Comparison at an arbitrary point within the support
- Exact solution vs gPC prediction for global radiation and precipitation
- $\triangleright$  For December 1970, large–scale patterns resolved in time-averaged results



Validation: Comparison of computed global means and gPC predictions

 $\triangleright$  Global mean should be consider to avoid small eccentric scales



## Sensitivity analysis

- $\blacktriangleright$  Partial variances reveal strong effects from 'entrpen'.
- $\triangleright$  Couple terms in the partial variance is much smaller
- Femperature PDF generated from the gPC metamodels with  $10^5$  Monte Carlo samples.



## Code calibration

For optimization problem with K objective functions, we seek all the  $\xi$  that satisfy the following minimization problem, e.g.

$$
\xi^* = \underset{\xi}{\text{argmin}} \sum_{k=1}^K \omega_k \left( \sum_{m=0}^M u_{m,k}(t) \phi_m(\xi) - u_{\text{obs},k}(t) \right)^2 \quad \text{for } t = 1,...,364
$$

The choice of weight vector  $\omega$  is arbitrary. Many optimization algorithms exist. So far  $K=1$ 

- $\blacktriangleright$  Lagrange multiplier algorithm used to solve the constraint nonlinear minimization problem for global averaged temperature
- $\triangleright$   $u_{\text{obs}}$  are the daily global averaged temperature in 1970 from ECMWF
- $\triangleright$  the following figures show the daily 'optimal' value for each parameter
- $\triangleright$  with additional objective functions, there is likely to be non-dominant sets, i.e. one cannot make one objective better without worsening the other objectives Neelin (2001)

## Calibration results





## Some concluding remarks

- $\triangleright$  PC and gPC constructs metamodels that accurately mimics the behaviours of complete simulators about the mean of the stochastic inputs
- Initial used as a UQ ans SA tool in engineering problems
- $\blacktriangleright$  It has potential as a multi-objective optimization tool
- $\triangleright$  There is no free lunch it suffers from the "curse of dimensionality"
- $\triangleright$  Adaptive techniques (multi-element, anisotropic quadrature) can reduce cost
- $\triangleright$  To investigate anisotropic spare quadrature & sparse gPC representation
- $\triangleright$  Reduce input dimension via non–dimensional analysis or identification of dominant inputs
- $\triangleright$  Orphan points (difference between sample budget and quadrature cost) can we use them in a sequential design – with Hugo?
- Including data assimilation and Bayesian analysis in  $gPC/PC$  framework
- $\blacktriangleright$  Practical issues: need better random input measurement