# Steps C: surrogate models

Michaël Baudin Chu Mai

May 20, 2021

#### Introduction

- ► The goal of this course is to introduce the main surrogate methods used in steps C.
- ▶ Principles of surrogate models, cross-validation, coefficient of predictability, over-fitting.
- ► Linear regression, polynomial chaos, kriging.
- ► The emphasis is not on maths.
- ▶ Instead, we feed the intuition with graphics.

References

Metamodels

Appendix

Some types of metamodels

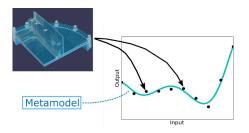
Methods for fitting metamodels

Training and validation

Some types of metamodels

Overfitting

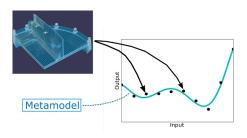
#### Metamodel - Definition



Meta: a prefix added to the name of something that consciously references or comments upon its own subject or features, e.g. metamodel: a model of another model

A metamodel is an approximation model that mimics the behavior of a computationally expensive simulator by training on *observations* (data) of the latter.

#### Metamodel - Definition



Expensive simulator:  $\mathbf{Y} = f(\mathbf{X})$ 

- $\triangleright$  X, Y are vectors of input and output,
- $X = (X_1, \dots, X_M)$

Metamodel:  $\tilde{Y} = \hat{f}(X, \theta)$ , where  $\theta$  is the vector of parameters By definition, two requirements of a metamodel  $\hat{f}$  are:

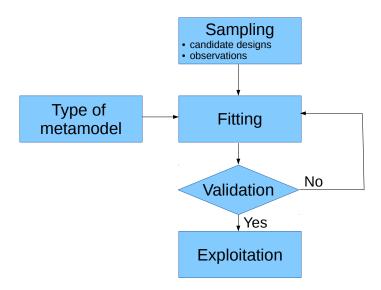
- ▶ Usefully accurate when predicting away from known observation
- ▶ Being significantly cheaper to evaluate than the primary simulator

M. Baudin

# Metamodel - Objectives

- ▶ Show functional relationships between input parameters and the output quantity of interest: impacts of variables
- ▶ Optimize the output quantity of interest: determine configurations that maximize the response or achieve specifications or customer requirements
- ► Replace the primary simulator in uncertainty propagation (surrogate model): sensitivity analysis

# Major steps for constructing a metamodel



#### Major steps for constructing a metamodel

#### Sampling (define an experimental design):

- ► A number of possible candidate designs are generated
- ▶ The designs are launched with the primary simulator

#### Constructing the metamodel:

- ► A type of metamodel is selected (among several available options)
- ▶ The metamodel is fitted to the available data
- ► The metamodel is validated (yes or no)
  - ▶ if yes: stop,
  - otherwise, reject the model.

# Major steps for constructing a metamodel

#### If the model is rejected:

- change method for fitting: e.g. use advanced regression technique instead of least squares errors,
- ▶ change metamodel parameters: e.g. increase polynomial degree,
- increase the sample size if the model is not accurate enough or interesting behavior is not observed,
- ► change the type of metamodel: e.g. polynomial chaos instead of second-order polynomial response surface

# Some types of metamodels

#### There are different metamodels available:

- ▶ Polynomial models
- ▶ Polynomial chaos models
- Kriging
- ► Neural networks,
- ▶ etc...

#### They all require to:

- ▶ assess the quality of the prediction: validation,
- $\triangleright$  tune the parameters  $\theta$ .

# Training and validation

Creating a metamodel involves two steps.

- 1. Train: estimate the coefficients of the metamodel which fits to the data,
- 2. Test: quantify the predictability of the metamodel.

One issue is to avoid overfitting, that is, we must be able to predict datasets that have not been used to train the metamodel (more on this topic later).

One solution is to use cross-validation.

- ▶ Split the sample into a train sub-sample and a test sub-sample,
- ► Train the metamodel on the train sample,
- ► Test the metamodel on the test sample.

#### Training and validation

Let  $\left\{\mathbf{x}_{t}^{(j)}\right\}_{j=1,\dots,n}$  an i.i.d. sample of the random vector  $\mathbf{X}$  that we use for training the metamodel.

We denote by g the model and  $\tilde{g}$  the metamodel. Let

$$y_t^{(j)} = g\left(\mathbf{x}_t^{(j)}\right), \qquad \tilde{y}_t^{(j)} = \tilde{g}\left(\mathbf{x}_t^{(j)}\right)$$

for j = 1, ..., n the outputs of the model and metamodel on the training set.

The  $R^2$  coefficient is:

$$R^{2}(g(\mathbf{x}_{t}), \tilde{g}(\mathbf{x}_{t})) = 1 - \frac{\sum_{j=1}^{N} \left(y_{t}^{(j)} - \tilde{y}_{t}^{(j)}\right)^{2}}{\sum_{j=1}^{N} \left(y_{t}^{(j)} - \bar{y}_{t}\right)^{2}}$$

where  $\bar{y}_t = \frac{1}{N} \sum_{i=1}^{N} y_t^{(j)}$ .

#### Training and validation

Denote  $\left\{\mathbf{x}_{v}^{(j)}\right\}_{j=1,\dots,n}$  the validation sample.

The goal of this sample is to test the metamodel on a dataset that was not used for training.

Let  $g(\mathbf{x}_v)$  and  $\tilde{g}(\mathbf{x}_v)$  the outputs of the model and metamodel on the validation sample.

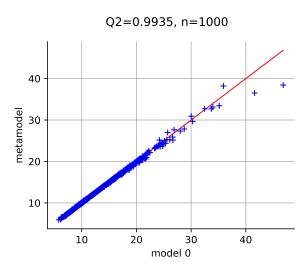
The  $Q^2$  predictability coefficient is defined as the  $\mathbb{R}^2$  applied to the validation sample:

$$Q^2 = R^2(g(\mathbf{x}_v), \tilde{g}(\mathbf{x}_v)).$$

A suggestion of decision rule:

- $ightharpoonup Q^2 > 0.95$ : accept the metamodel,
- otherwise: try to improve the metamodel or reject it.

# Graphical validation



# Overfitting

Overfitting is the fact that the surrogate model fits to the training sample, but predicts poorly on new points.

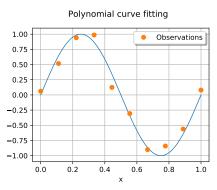
- ► Any surrogate has coefficients tuned on the training sample.
- ▶ More coefficients generally reduce the error on the training sample
- but may increase the error on the validation sample.

There is a bias-variance trade-off to solve:

- ► More coefficients may reduce the variance (seen on the training sample),
- but may increase the bias (revealed on the validation sample).

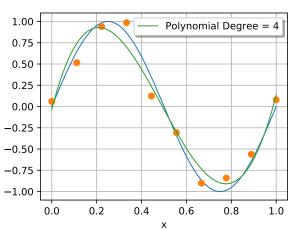
#### Example 1

- ▶ We have 10 points on the [0,1] interval, produced by adding a small noise to the sine function.
- ► We approximate it with linear regression based on polynomial canonical basis.

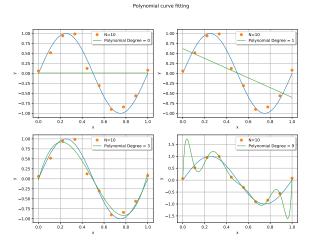


We use linear least squares to fit a degree 4 polynomial (5 coefficients).

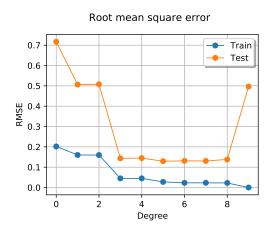




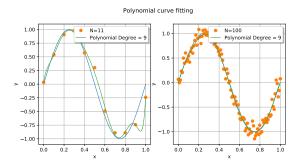
Increasing the degree may not always reduce the generalization error.



When we increase the degree, the root mean square error (RMSE) first decreases, then increases.



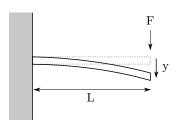
Adding points to the training sample reduces the validation error, ...



 $\dots$  but this is not always possible, e.g. the computer code may require costly simulations.

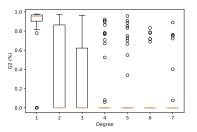
#### Example 2

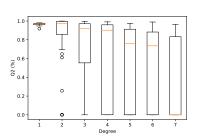
- Consider the cantilever beam example  $Y = \frac{FL^3}{3EI}$  with independent marginals: E: Beta, F: Lognormal, L: Uniform, I: Beta.
- ▶ We approximate it with *sparse* polynomial chaos: this limits the number of coefficients in the model.

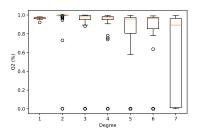


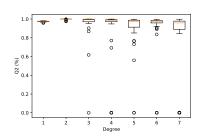
#### Questions:

- $\blacktriangleright$  How to set the sample size n?
- $\blacktriangleright$  How to set the polynomial degree d?







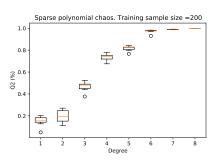


#### Answer:

- ▶ How to set the sample size n? n > 40
- ▶ How to set the polynomial degree d? d = 2 is the smallest degree with high  $Q_2$

#### Example 3

- Consider the Ishigami model  $Y = \sin(X_1) + a\sin(X_2)^2 + bX_3^4\sin(X_1)$  where a = 7 and b = 0.1, with independent marginals:  $X_1, X_2, X_3 \sim \mathcal{U}(-\pi, \pi)$ .
- ▶ We approximate it with *sparse* polynomial chaos: this limits the number of coefficients in the model.



#### Références I



Christopher M Bishop et al., Neural networks for pattern recognition, Oxford university press, 1995.

Appendix

Appendix

#### Some types of metamodels

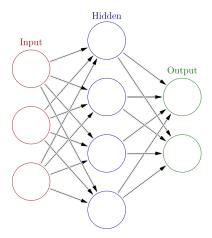
#### Radial basis function:

$$\tilde{\boldsymbol{Y}} = \sum_{k=1}^{N} \theta_k \psi(\|\boldsymbol{X} - \boldsymbol{X}_k\|)$$

where  $\psi()$  being a radial basis function with its centers taken at  $X_k$ , k = 1, ..., N in the experimental design

#### Some types of metamodels

**Artificial neural network:** Radial basis function is a single layer neural network with radial coordinate neurons



Regularized regression methods: minimize sum of squared errors under a constraint

$$\boldsymbol{\theta}^* = \arg\min_{\boldsymbol{\theta}} \sum_{i=1}^{N} (Y_i - \hat{f}(\boldsymbol{X}_i, \boldsymbol{\theta}))^2 + \lambda R(\boldsymbol{\theta})$$

- $ightharpoonup R(\theta) = ||\theta||_2$ : Ridge regression,
- $ightharpoonup R(\theta) = ||\theta||_1$ : LASSO regression,
- $\triangleright$   $\lambda$ : non-negative regularization coefficient

# Methods for fitting metamodels

Least square regression: minimize sum of squared errors of a linear regression model

$$\boldsymbol{\theta}^* = \arg\min_{\boldsymbol{\theta}} \sum_{i=1}^N \epsilon_i^2 = \arg\min_{\boldsymbol{\theta}} \sum_{i=1}^N (Y_i - \hat{f}(\boldsymbol{X}_i, \boldsymbol{\theta}))^2$$

#### Methods for fitting metamodels

Maximum likelihood estimation: e.g. assume that the errors  $\epsilon$  are independently randomly distributed according to a normal distribution with standard deviation  $\sigma$ 

$$\mathcal{L}(\boldsymbol{\theta}) = \frac{1}{(2\pi\sigma^2)^{N/2}} \prod_{i=1}^{N} \exp\left(-\frac{1}{2} \left(\frac{Y_i - \hat{f}(\boldsymbol{X}_i, \boldsymbol{\theta})}{\sigma}\right)^2\right)$$
$$\boldsymbol{\theta}^* = \arg\max_{\boldsymbol{\theta}} \mathcal{L}(\boldsymbol{\theta})$$

# Methods for fitting metamodels

#### *K*-fold cross-validation method:

 $\mathcal{K}: \{1,\ldots,N\} \mapsto \{1,\ldots,K\}$  partition of N observations to K roughly equal-sized parts, K=N: leave-one-out  $\hat{f}^{-k}()$ : fitted metamodel with k-th part of data set aside Cross-validation estimate of prediction error:

$$CV(\hat{f}, \boldsymbol{\theta}) = \frac{1}{N} \sum_{i=1}^{N} L(Y^i, \hat{f}^{-\mathcal{K}(i)}(X_i))$$
$$\boldsymbol{\theta}^* = \arg\min_{\boldsymbol{\theta}} CV(\hat{f}, \boldsymbol{\theta})$$

# Some types of metamodels

Polynomial models: (response surface)

Second-order model

$$\tilde{Y} = \theta_0 + \sum_{i=0}^{M} \theta_i X_i + \sum_{i=0}^{M} \theta_{ii} X_i^2 + \sum_{i < j} \sum_{j=2}^{M} \theta_{ij} X_i X_j$$

#### Polynomial chaos expansion:

$$\tilde{\boldsymbol{Y}} = \sum_{0 \leq |\boldsymbol{k}| \leq p} \, \theta_{\boldsymbol{k}} \psi_{\boldsymbol{k}}(\boldsymbol{X})$$

where  $\psi_{\mathbf{k}}()$  being polynomial chaos functions

#### Some types of metamodels

**Kriging:** (Gaussian process regression)

Deterministic trend: linear regression on a fixed basis

$$m(\boldsymbol{X}) = \boldsymbol{r}(\boldsymbol{X})^T \boldsymbol{\theta}$$

Random fluctuation: zero-mean stationary Gaussian process of covariance function

$$k(\boldsymbol{X}, \boldsymbol{X'}) = \sigma^2 \rho(\|\boldsymbol{X} - \boldsymbol{X'}\|)$$