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SPECIAL MOBILITY STRAND

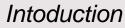
Reliability Analysis Techniques Juan Sepúlveda Novi Sad, Serbia, 22nd February 2020

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Presentation outline





Deterministic design Probabilistic design Variable transformation

First Order Reliability Method (FORM)

Second Order Reliability Method (SORM)

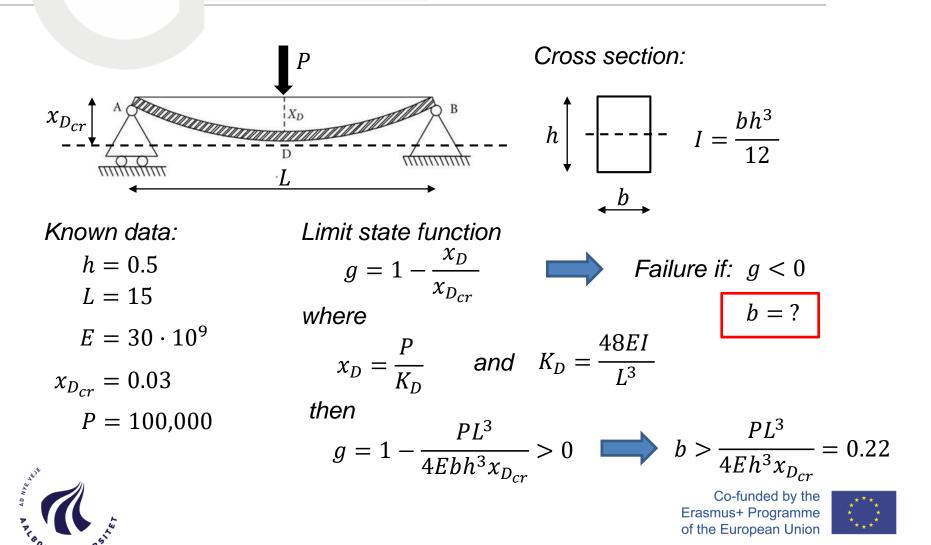
Crude Monte Carlo Simulation (CMCS)

Importance Sampling (IS)

Asymptotic Samplig (AS)



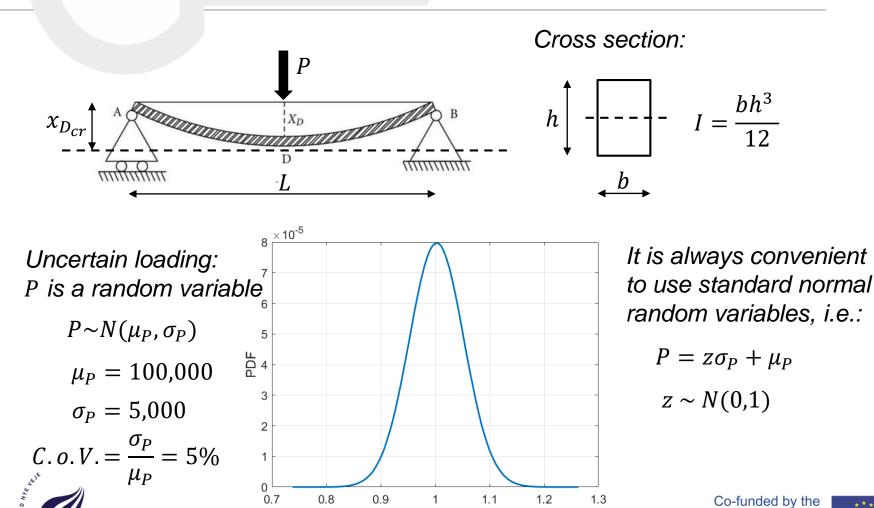






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 $\times 10^5$



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A simple transformation from X_i to Z_i can be defined by the identity

 $\Phi(Z_i) = F_{X_i}(X_i)$

where F_{X_i} is the distribution function for X_i . Given a realization z of Z a realization x of X can be determined by

 $x_i = F_{X_i}^{-1}(\Phi(z_i))$





The failure probability is given as

$$P_F = P(g < 0) = \int_F q(\mathbf{z}) \mathrm{d}\mathbf{z}$$

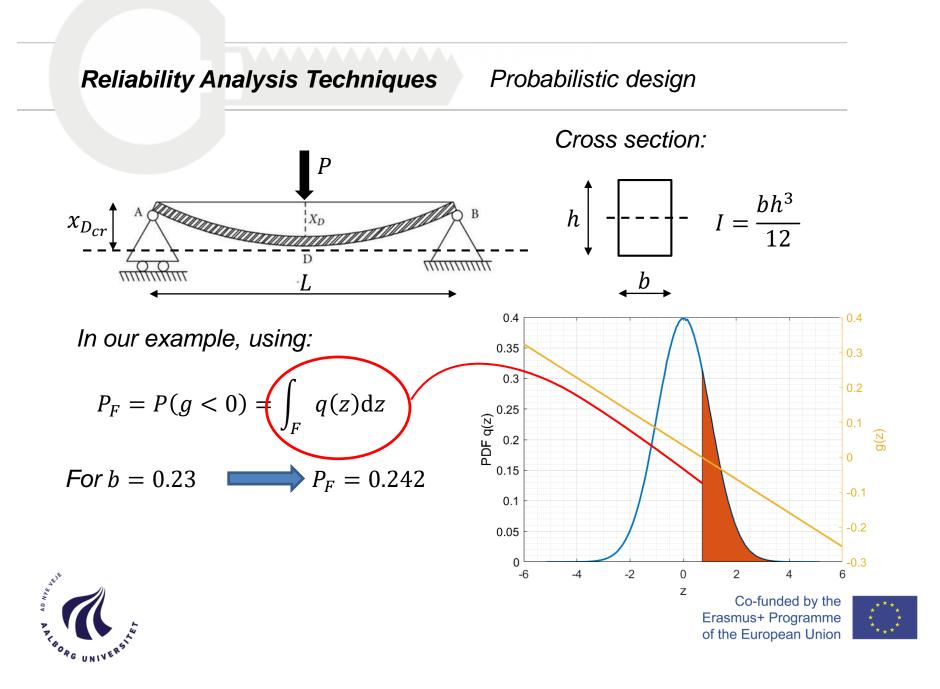
Where $q(\mathbf{z})$ is the joint Probability Density Function (PDF) of the random variables and

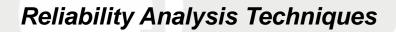
$$F = \{ \mathbf{z} : g < 0 \}$$

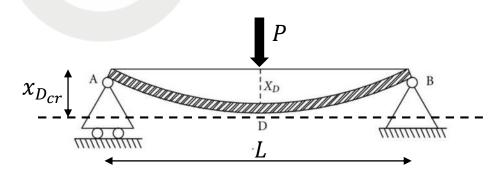
Denotes the "failure domain" or "failure region" which is a subset in the parameter space of z



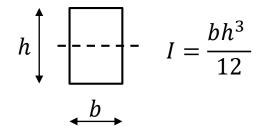








Cross section:

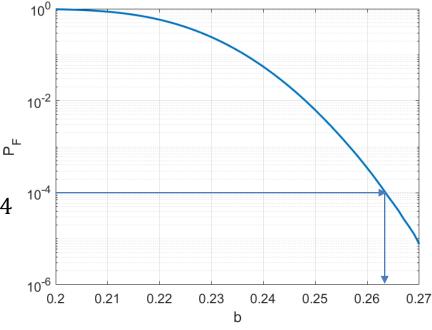


The design requirement now is the target reliability of the system

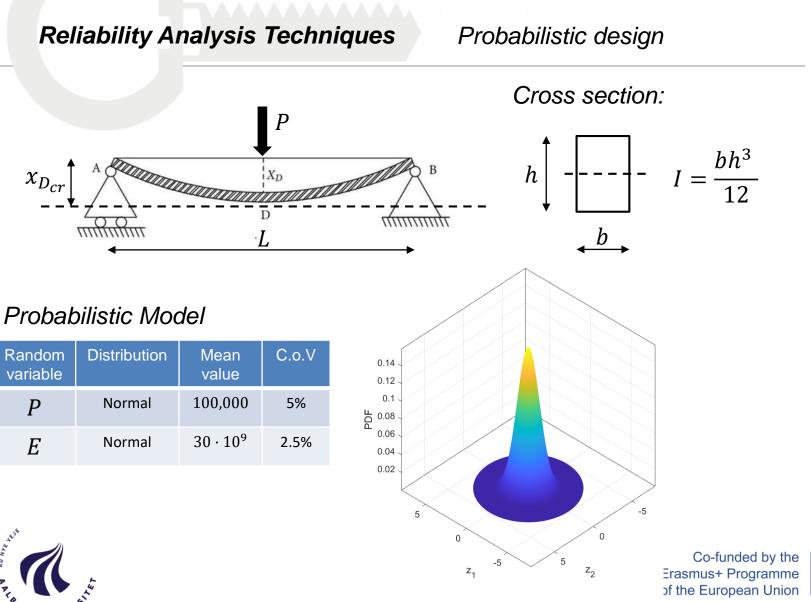
 $R^* = 1 - P_F^*$

Or equivalently the target failure probability P_F^*

For example: $P_F^* = 10^{-4}$ $\implies b = 0.264$





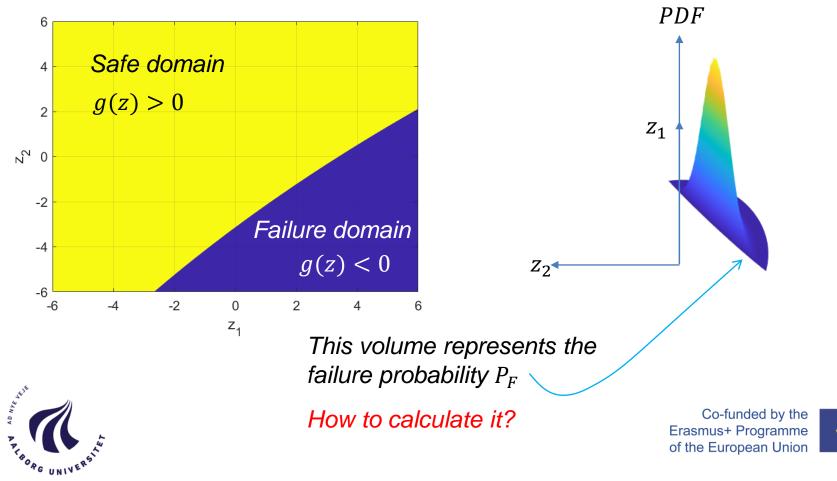


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Probabilistic design

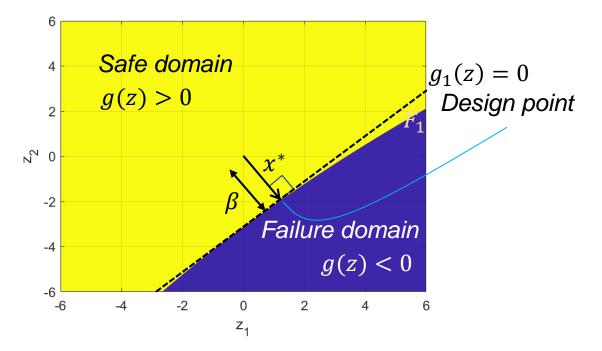
For a given value of b = 0.26



The failure domain is approximated by a linear half-space F_1 .

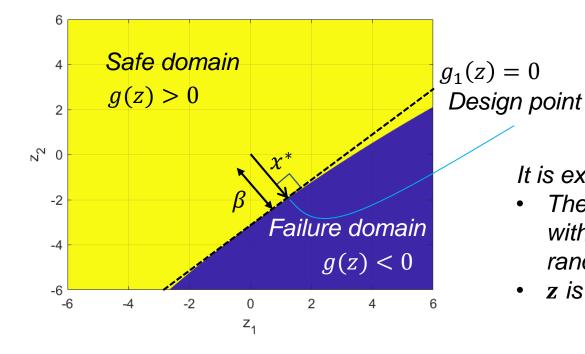
The "design point" is the point in F_1 with minimal distance to the origin, i.e. with minimal Euclidean norm.

The maximum of the PDF in F_1 is at the design point.









The approximation is given by $P_F \approx \widehat{P_F} = \Phi(-\beta)$

It is exact when:

- The limit state function is linear with respect to the vector of random variables z
- z is Gaussian distributed



For b = 0.26, the approximation gives $P_F \approx \widehat{P_F} = \Phi(-2.243) = 0.0124$



How to determinate β ?

- 1. Guess $z^{(0)}$. Set i = 0.
- 2. Calculate $g(\mathbf{z}^{(i)})$ and $\nabla g(\mathbf{z}^{(i)})$
- 3. Calculate an improved guess of β with $\mathbf{z}^{(i+1)} = \nabla g(\mathbf{z}^{(i)}) \frac{\nabla g(\mathbf{z}^{(i)})^T \mathbf{z}^{(i)} - g(\mathbf{z}^{(i)})}{\nabla g(\mathbf{z}^{(i)})^T \nabla g(\mathbf{z}^{(i)})}$
- 4. Calculate

$$\beta^{(i+1)} = \sqrt{(\mathbf{z}^{(i+1)})^T \mathbf{z}^{(i+1)}}$$

5. If $|\beta^{(i+1)} - \beta^{(i)}| < 10^{-3}$ then stop. Else i = i + 1 and go to step 2. Co-funded by the Erasmus+ Programme of the European Union



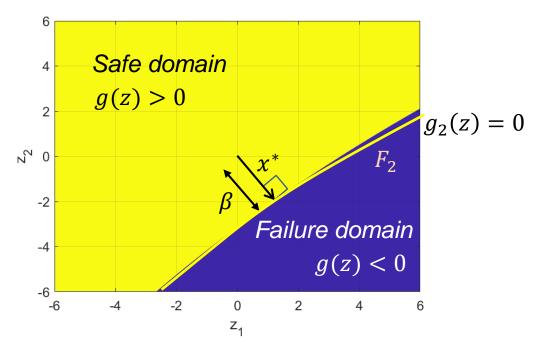
- The estimator does not depend of the order of $\widehat{P_F}$
- Under particular conditions leads to the exact solution
- The linearization introduce an error
- It is found that is not accurate in high dimensions
- It is necessary to find the design point
 - It might be expensive in high dimensions





It intends to improve FORM taking into account the curvature of limit state function at the design point.

The failure boundary is approximated by an hyperparaboloid with limit state function g_2 .







The SORM approximation is given by

$$P_F \approx \widehat{P_F} = \Phi(-\beta) \prod_{i=2}^n (1+c_i\beta)^{-\frac{1}{2}}, \qquad \beta \to \infty$$

Where c_i , i = 2, ..., n are the principal curvatures of the paraboloid at the design point.

It is also assumed that $-c_i < 1/\beta$ for all i = 2, ..., n



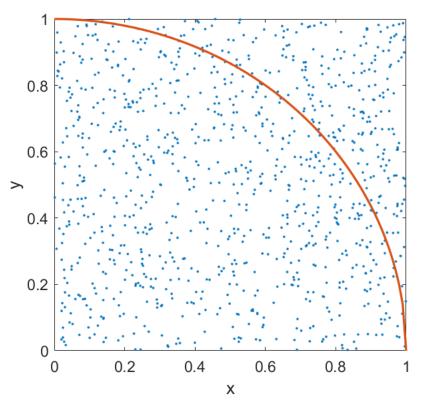


How to calculate the area of a circle? And implicitly π ?

- Simulate N random numbers into a unit square with area A = 1
- Draw a $\frac{1}{4}$ of a unit-radio circle with area $A_1 = \pi/4$
- Count the points inside the circle N_1

• The ratio
$$\frac{A_1}{A} = \frac{N_1}{N}$$

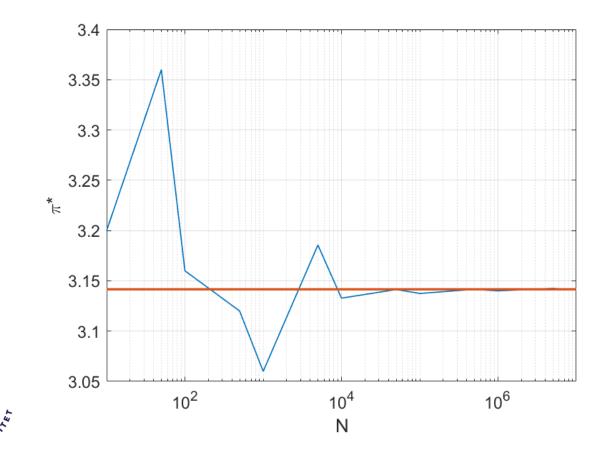






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How is affected the estimation of $\pi \approx \pi^*$ by the number of simulations?

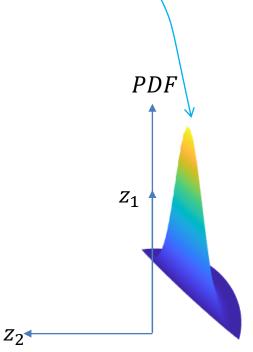






We can use the same approach to calculate this volume -

- A large number N of realizations ẑ of basic random variables is simulated using q(z)
- The failure probability estimator $\widehat{P_F}$ is obtained by counting the number of realizations in the failure domain



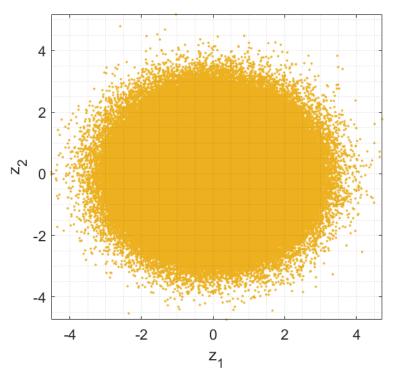




Crude Monte Carlo Simulation

Procedure

- N realizations of standard normal random variables are simulated
 - This is the only step where the information regarding the distribution of the random variables in included





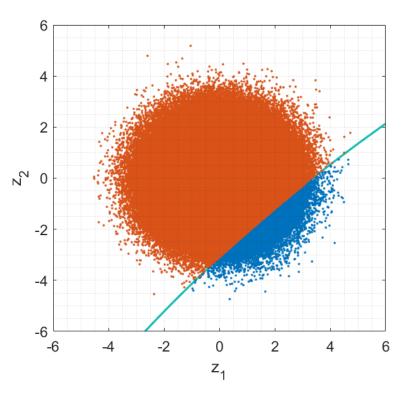




Crude Monte Carlo Simulation

Procedure

- *1. N* realizations of standard normal random variables are simulated
 - This is the only step where the information regarding the distribution of the random variables is included
- 2. Identify the number N_1 of realizations in the failure domain
 - This is the only step where the information regarding the limit state function is included





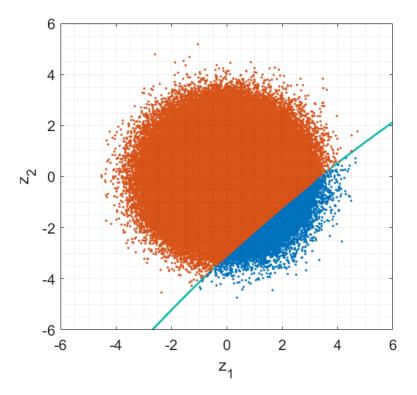


Procedure

- *1. N* realizations of standard normal random variables are simulated
 - This is the only step where the information regarding the distribution of the random variables is included
- 2. Identify the number N_1 of realizations in the failure domain
 - This is the only step where the information regarding the limit state function is included
- 3. The estimator is given by

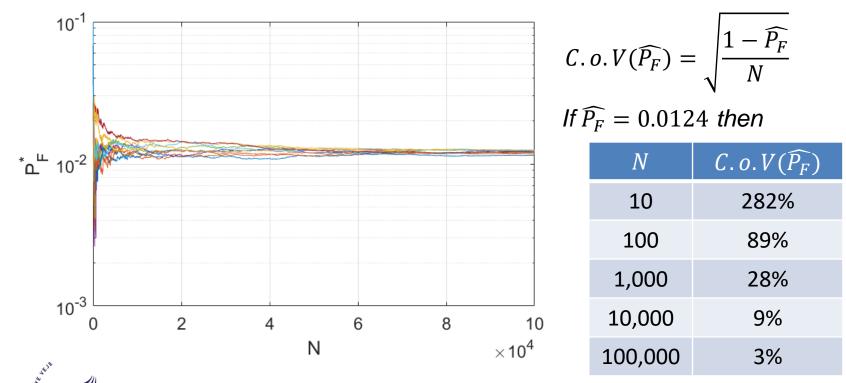






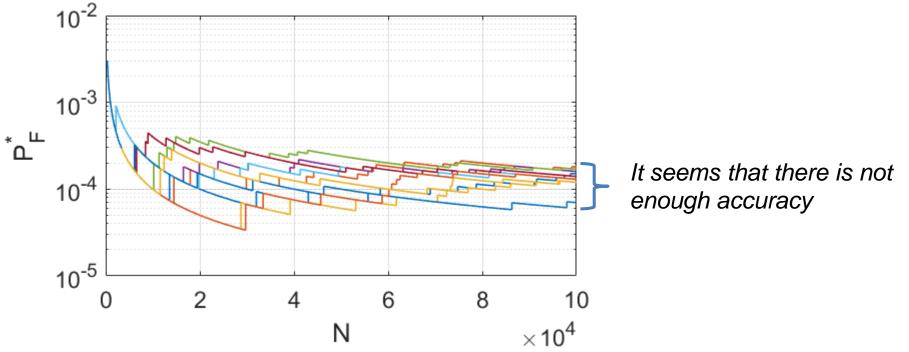


In our example, using b = 0.26, how N affects the variability of the estimator $\widehat{P_F}$?





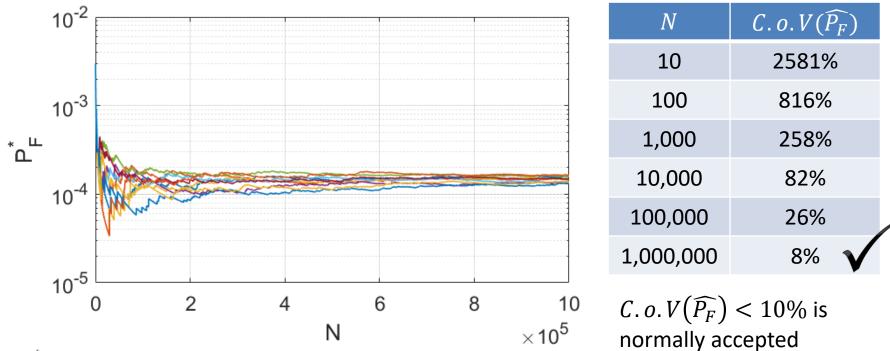
In our example, using b = 0.285, i.e. $\widehat{P_F} = 1.5 \times 10^{-4}$. How $\widehat{P_F}$ affects C.o.V($\widehat{P_F}$)?





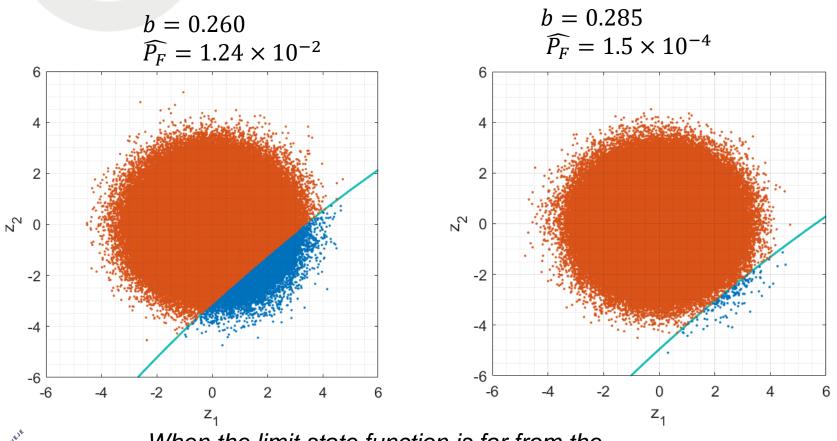


In our example, using b = 0.285, i.e. $\widehat{P_F} = 1.5 \times 10^{-4}$. How $\widehat{P_F}$ affects $C.o.V(\widehat{P_F})$?



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When the limit state function is far from the origin is more difficult to obtain realizations in the failure domain



Crude Monte Carlo Simulation



- Independent of the number of random variables
- Independent of the shape of the limit state function
- The estimator is unbiased

• The number of samples required is proportional to $1/P_F$





Importance Sampling

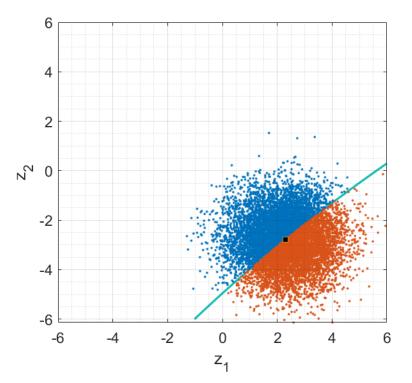
The idea is to concentrate the realizations in an area with large contribution to $\widehat{P_F}$

$$P_F = \int_{g<0} q(\mathbf{z}) d\mathbf{z} = \int_{g<0} \frac{q(\mathbf{z})}{p(\mathbf{z})} p(\mathbf{z}) d\mathbf{z}$$

Then, the estimator is given by

$$\widehat{P_F} = \frac{1}{N} \sum_{k=1}^{N} \frac{q(\mathbf{z}'_k)}{p(\mathbf{z}'_k)} I(\mathbf{z}'_k \in F)$$

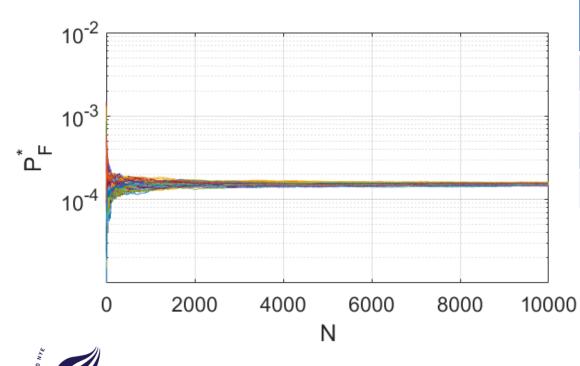
Where $\{\mathbf{z}'_k: k = 1, ..., N\}$ are now samples drawn from p instead of q







In our example, using b = 0.285, i.e. $\widehat{P_F} = 1.5 \times 10^{-4}$. How N affects the variability of the estimator $\widehat{P_F}$?



N	$C.o.V(\widehat{P_F})$ MCS	$C.o.V(\widehat{P_F})$ IS
10	2581%	65%
100	816%	19%
1,000	258%	7% 🗸
10,000	82%	2% 🗸

With 1,000 samples achieves the same accuracy than 1,000,000 samples using CMCS



Importance Sampling



- Reduce dramatically the number of samples required by using CMCS
- An incorrect selection of the importance sampling density function can lead to erroneous estimates of the reliability





The idea is to use asymptotic properties with a simple regression technique

- CMCS is used to estimate the failure probability in each support point
- The standard deviation for each support point is increased by a factor $\frac{1}{f}$
- Due to the larger standard deviation the failure probability is increased reducing the computational cost
- The functional form is given by

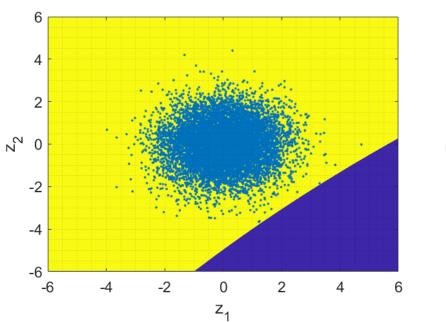
$$\beta(f) = A \cdot f + \frac{B}{f^c}$$

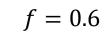


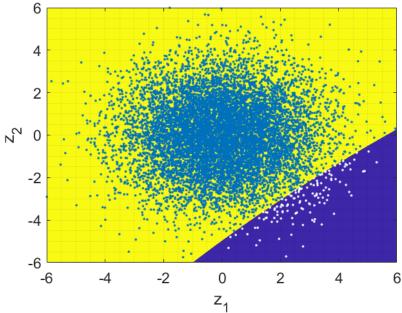


f = 1

Asymptotic Sampling

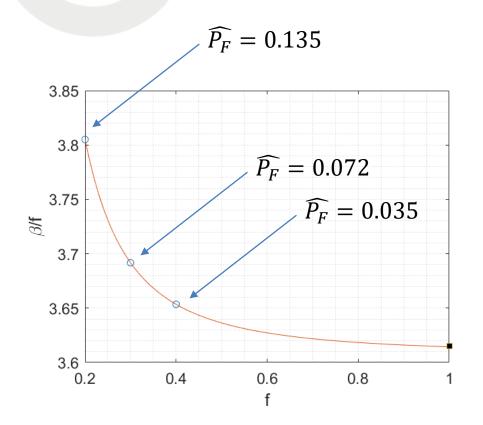












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The failure probabilities are much larger than the original implying a lower computational cost

Asymptotic Sampling

Performing a regression, the following parameters are estimated

- A = 3.6078B = 0.0066
- C = 2.1087

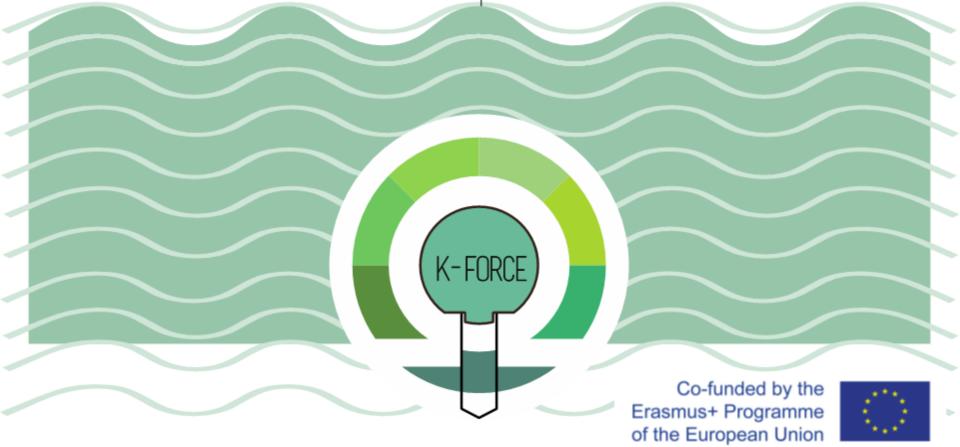
Then the functional form

 $\beta(f) = A \cdot f + \frac{B}{f^{C}}$ Is evaluated for f=1 $\beta(1) = 3.61$

And then the estimator is given by

$$\widehat{P_F} = \Phi(-\beta(1)) = 1.5 \times 10^{-4}$$





Thank you for your attention jgsa@civil.aau.dk

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