

Response Surface Methodology: Central Composite Designs

BIOE 498/598 PJ

Spring 2021

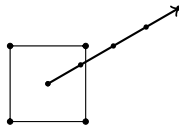
Last time: The method of steepest ascent

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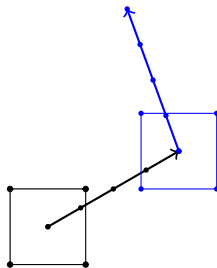
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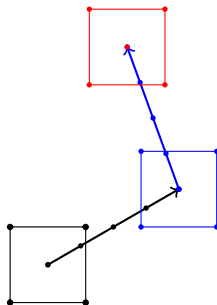
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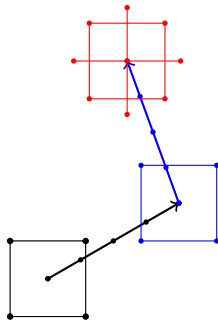
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- ▶ Stop when model detects lack of fit.
- ▶ **Today:** Fitting a model to a curved response surface.



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$$y = 20 + 3.6x_1 - 1.8x_2 - 0.6x_1x_2$$

Set $x_2 = 0$, then $y \rightarrow \infty$ as $x_1 \rightarrow \infty$.

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- ▶ If you know f for your system, congrats! Fit its parameters with regression and use it.
- ▶ Usually we don't know f , so we approximate it with a simpler function.
- ▶ **We are not claiming that f is a particular shape.** Rather, we claim that an approximation is “good enough” over our domain of interest.

Approximating f with a general quadratic

Let's find the second-order Taylor series of $f(x_1, x_2)$ centered at zero:

$$f(x_1, x_2) \approx \underbrace{f|_0}_{\beta_0} + \underbrace{\frac{\partial f}{\partial x_1} \Big|_0}_{\beta_1} x_1 + \underbrace{\frac{\partial f}{\partial x_2} \Big|_0}_{\beta_2} x_2 + \underbrace{\frac{1}{2} \frac{\partial^2 f}{\partial x_1^2} \Big|_0}_{\beta_{11}} x_1^2 + \underbrace{\frac{1}{2} \frac{\partial^2 f}{\partial x_2^2} \Big|_0}_{\beta_{22}} x_2^2 + \underbrace{\frac{1}{2} \frac{\partial^2 f}{\partial x_1 \partial x_2} \Big|_0}_{\beta_{12}} x_1 x_2$$

$$f(x_1, x_2) \approx \underbrace{\beta_0 + \beta_1 x_1 + \beta_2 x_2}_{\text{FO}} + \underbrace{\beta_{11} x_1^2 + \beta_{22} x_2^2}_{\text{PQ}} + \underbrace{\beta_{12} x_1 x_2}_{\text{TWI}}$$

SO

- ▶ The function f and its derivatives are unknown, so we fit the parameters β with a linear model.
- ▶ In general we will have k factors and the quadratic approximation will be

$$y = \beta_0 + \sum_{i=1}^k \beta_i x_i + \sum_{i=1}^k \beta_{ii} x_i^2 + \sum_{j=1}^k \sum_{i=1}^{j-1} \beta_{ij} x_i x_j.$$

- ▶ This model has $1 + 2k + k(k-1)/2$ parameters, so RSM designs must have at least this many runs.

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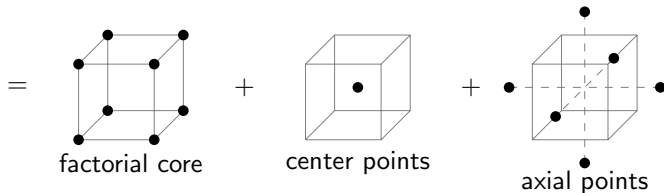
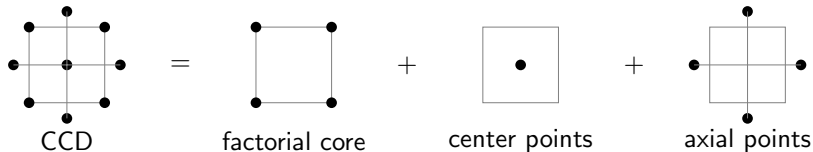
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The Central Composite Design (CCD)

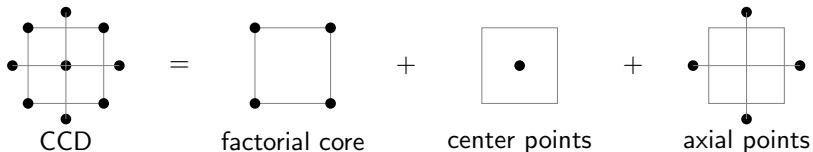
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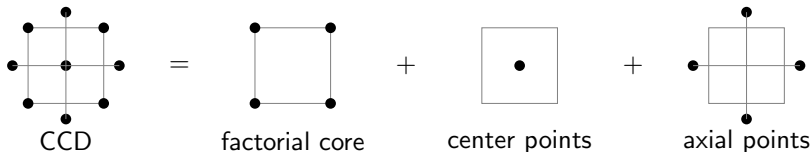


Parts of the CCD



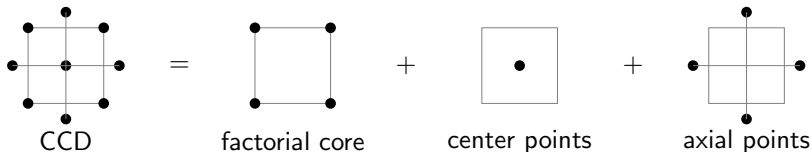
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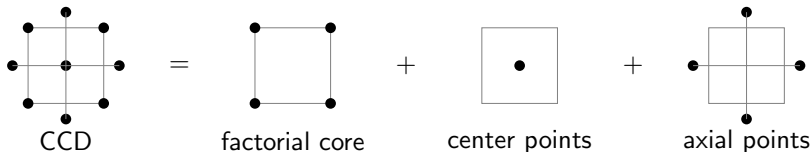
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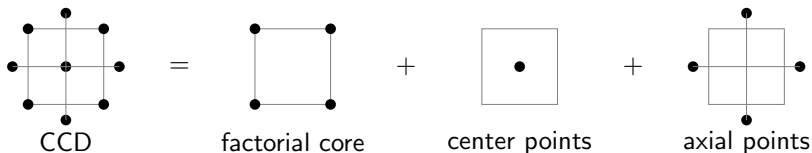
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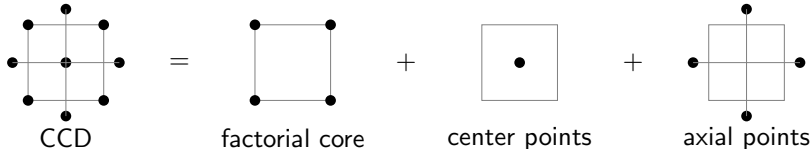
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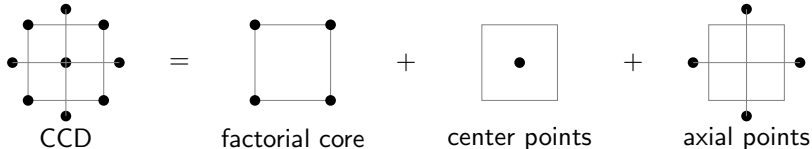
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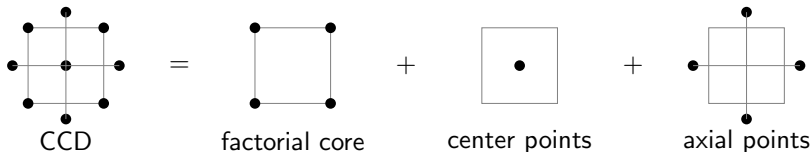
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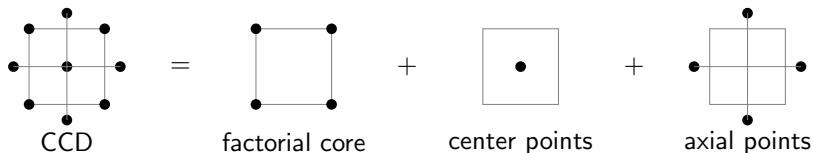
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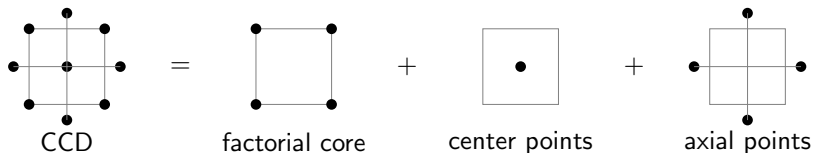
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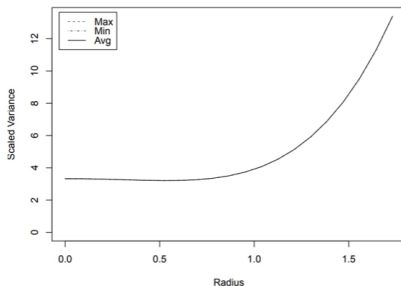
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 3. The value of α

Uniform precision

- ▶ A model has **uniform precision** if the variance at design radius 1 is the same as at the center.

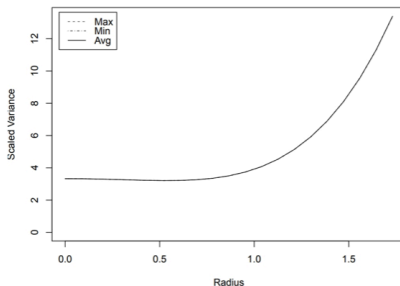
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- ▶ Choosing the correct number of center points in a CCD ensures uniform precision.

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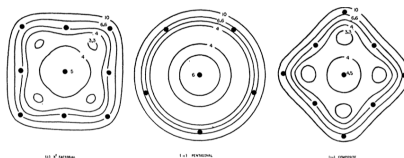


FIG. 2. Variance contours for some 2 dimensional designs

Image from Box and Hunter 1957.

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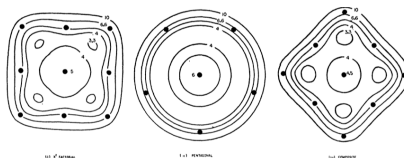


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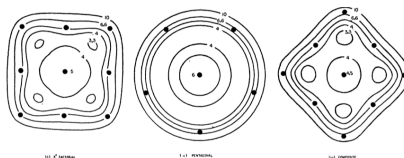


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- ▶ A CCD with F factorial points is rotatable when $\alpha = \sqrt[4]{F}$.

Rotatable, uniform precision CCDs

factors (k)	2	3	4	5	5 - 1	6
factorial points	4	8	16	32	16	64
axial points	4	6	8	10	10	12
center points	5	6	7	10	6	15
axial distance (α)	1.414	1.682	2.000	2.378	2.000	2.828

factors (k)	6 - 1	7	7 - 1	8	8 - 1	8 - 2
factorial points	32	128	64	256	128	64
axial points	12	14	14	16	16	16
center points	9	21	14	28	20	13
axial distance (α)	2.378	3.364	2.828	4.000	3.364	2.828

Factor levels in a CCD

Each factor in the CCD will be set at five levels:

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Unlike a 2-level design, the coded units in a CCD have meaning!

Coding the CCD

Let's say we're designing a combination screening of three drugs. The absolute widest concentration range we can use for drug A is $[-3.2, 1.0]$ on a \log_{10} - μM scale. What are the five levels assuming a full-factorial CCD?

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$$\begin{aligned} A &= \text{center}(A) + \frac{\text{range}(A)}{2\alpha}[\text{code}] \\ &= -1.1 + \frac{1 - (-3.2)}{2(1.68)}[\text{code}] \end{aligned}$$

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Let's say we're designing a combination screening of three drugs. The absolute widest concentration range we can use for drug A is $[-3.2, 1.0]$ on a \log_{10} - μ M scale. What are the five levels assuming a full-factorial CCD?

$$F = 2^3 = 8 \Rightarrow \alpha = \sqrt[4]{8} = 1.68$$

$$\begin{aligned} A &= \text{center}(A) + \frac{\text{range}(A)}{2\alpha}[\text{code}] \\ &= -1.1 + \frac{1 - (-3.2)}{2(1.68)}[\text{code}] \end{aligned}$$

code:	$-\alpha$	-1	0	1	α
\log_{10} - μ M:	-3.2	-2.4	-1.1	0.2	1.0