



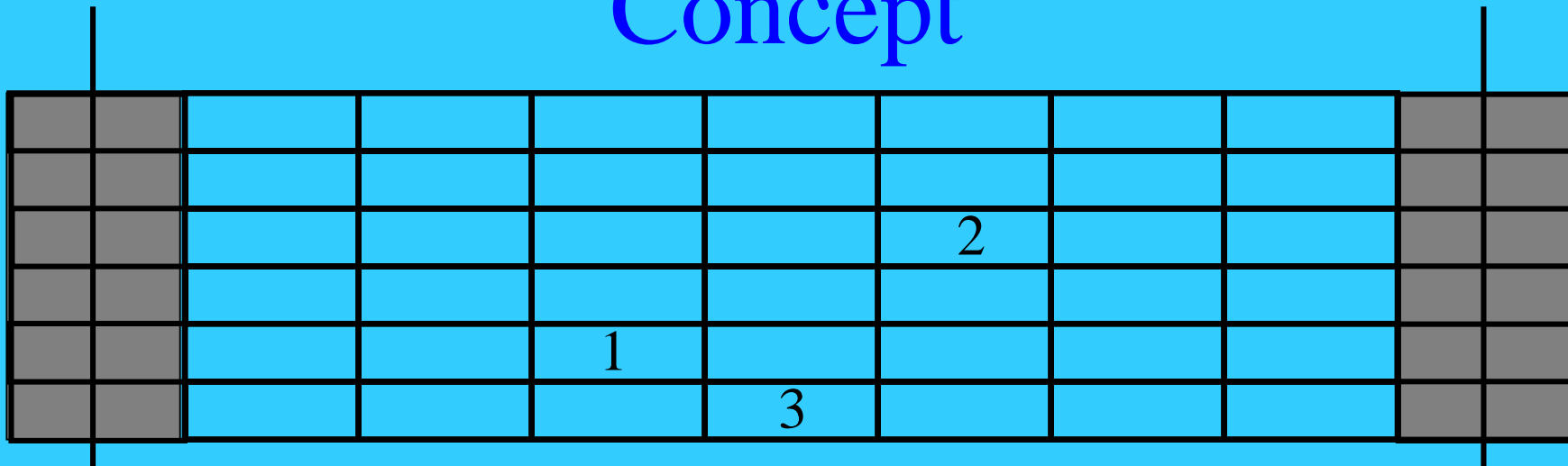
Reservoir Modeling with GSLIB

# Indicator Simulation for Categorical Variables

- Sequential Simulation: the Concept
- Steps in Sequential Simulation
- SISIM Program



# Sequential Simulation: the Concept

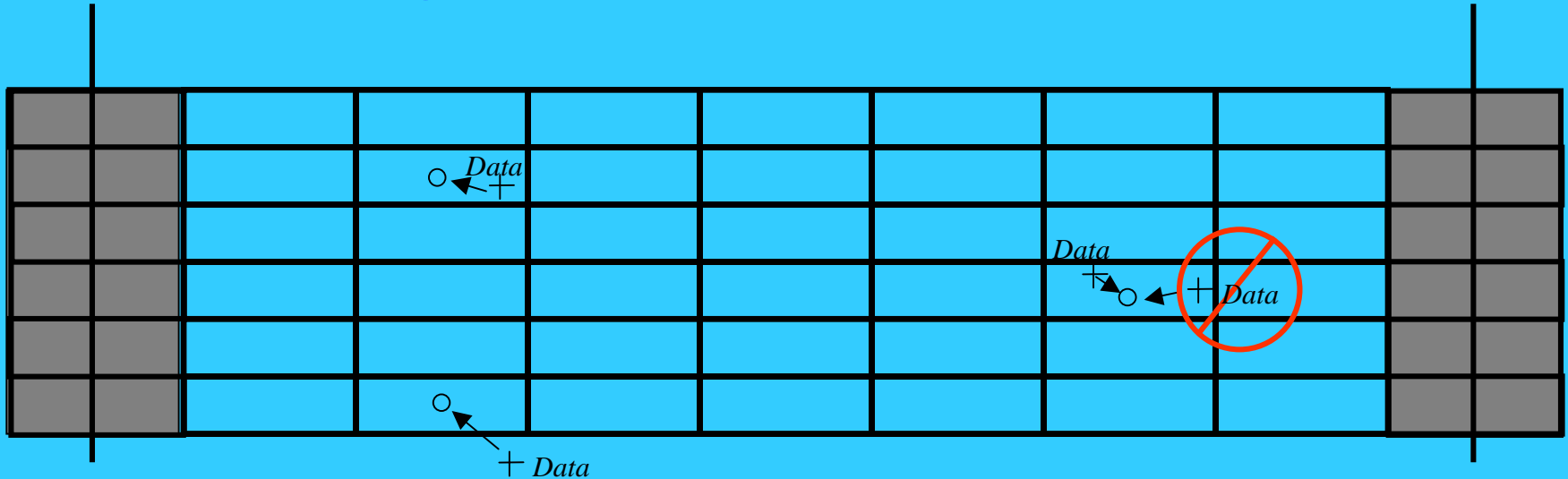


1. Assign data values to closest grid node
2. Establish a random path through all of the grid nodes
3. Visit each grid node:
  - (a) find nearby data and previously simulated grid nodes
  - (b) construct the conditional distribution by kriging (this is where the variogram comes in)
  - (c) draw simulated value from conditional distribution
4. Check the results



## STEP 1


# Assign Data to Grid Nodes



Why?

- Explicitly honor data  $\Rightarrow$  data values will appear in final 3-D model
- Improves the CPU speed of the algorithm: searching for previously simulated nodes and original data is accomplished in one step

Considerations:

- Take the closest of multiple data within the same cell  $\Rightarrow$  could lose some information (see in figure )
- Not an option when simulating a cross-sectional or small-area model



## STEP 2

# Establish a Random Path

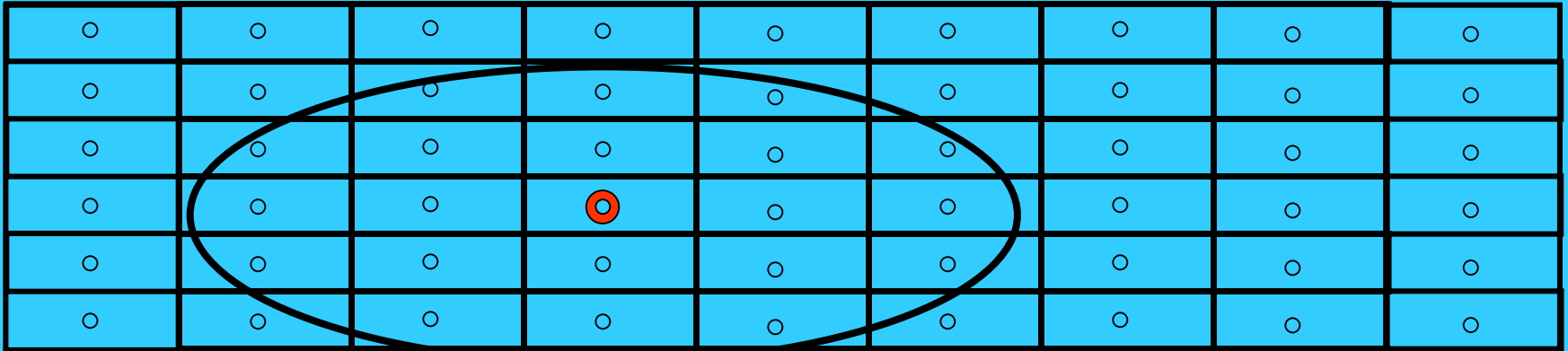
|  |  |    |    |    |    |    |    |    |  |  |
|--|--|----|----|----|----|----|----|----|--|--|
|  |  | 32 | 4  | 34 | 18 | 35 | 2  | 21 |  |  |
|  |  | 30 | 31 | 14 | 8  | 9  | 16 | 3  |  |  |
|  |  | 28 | 6  | 24 | 5  | 15 | 13 | 18 |  |  |
|  |  | 11 | 5  | 28 | 20 | 27 | 23 | 10 |  |  |
|  |  | 1  | 27 | 19 | 33 | 25 | 4  | 26 |  |  |
|  |  | 29 | 29 | 7  | 12 | 17 | 6  | 22 |  |  |

- Visit each cell once and only once in random order
- Can do this in many ways:
  - draw a random number and multiply it by  $N$
  - sort an array of random numbers while carrying an array of indices  
capitalize on the limited period length of linear congruential generators
- Skip over cells (actually grid nodes) that already have a value



## STEP 3 - (a)

# Find Nearby “Informed” Nodes

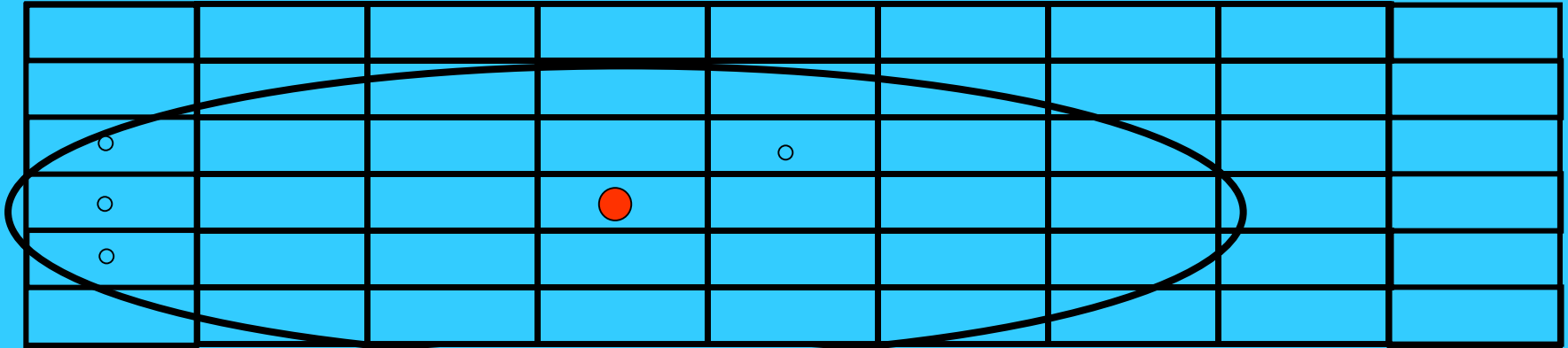


- “Informed” nodes refers to both data-nodes and nodes that have been informed earlier in the random path
- Typically use spiral search to identify the close nodes
- Limit the number of nodes actually considered:
  - octant search (?)
  - maximum per octant (say 4)
  - maximum number



STEP 3 - (b)

# Construct Conditional Distribution



- Conditional distribution is constrained by:
  - global proportion of each lithology type
  - local data
  - “local” proportion from secondary data such as seismic
- Calculate by kriging the binary indicator transform for each rock type



# Indicator Simulation (1)

- Define an indicator transform:

$$i(u_{\alpha}; k) = \left\{ \begin{array}{l} 1, \text{ if lithofacies } k \text{ present at location } u_{\alpha} \\ 0, \text{ if not} \end{array} \right\}$$

- Average of an indicator is the global proportion:

$$\text{proportion of } k = E\{I(u_{\alpha}; k)\}$$

$$= \frac{\sum_{\alpha=1}^n w_{\alpha} \cdot i(u_{\alpha}; k)}{n}$$

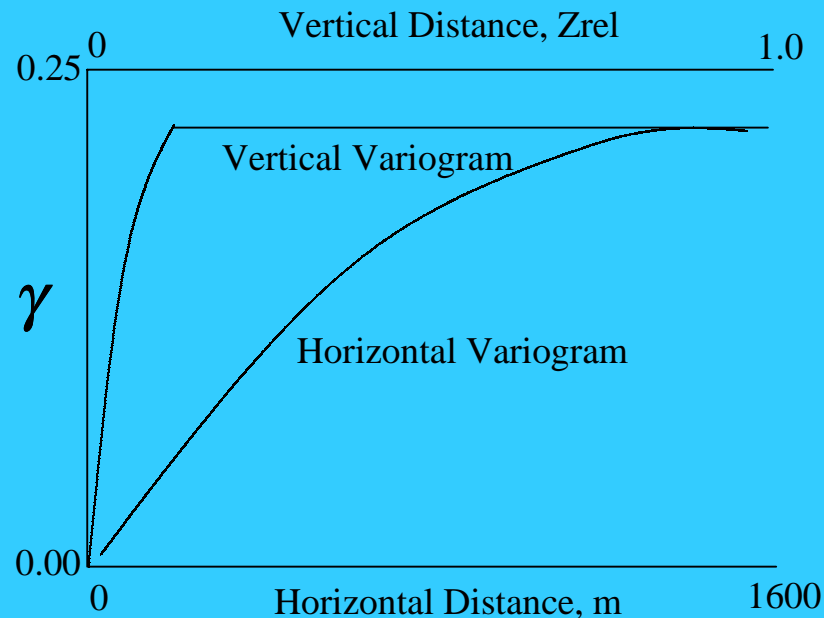
$w_{\alpha}$  values account for data clustering



# Indicator Simulation (2)

- The variogram of an indicator variable measures spatial correlation:

$$\gamma_I(h) = \frac{1}{2} E \{ [I(u; k) - I(u + h; k)]^2 \}$$

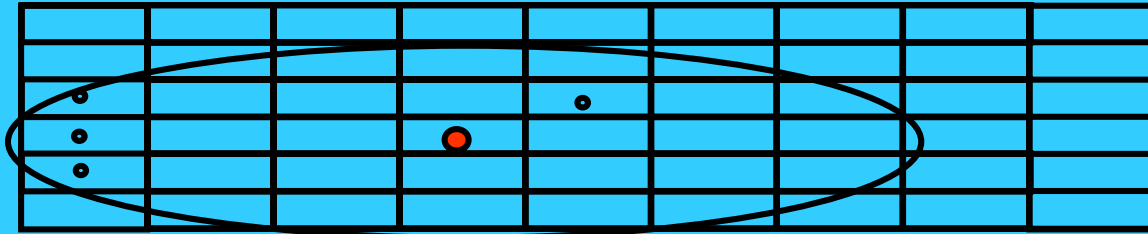






STEP 3 - (c)

# Construct Conditional Distribution with Kriging



- Given  $n$  nearby data values  $k(\mathbf{u}_i), i=1, \dots, n$  how do we calculate the conditional distribution?
- Estimate conditional probabilities of each rock type  $p_k^*(\mathbf{u}), k=1, \dots, K$  by a linear combination of the nearby data:

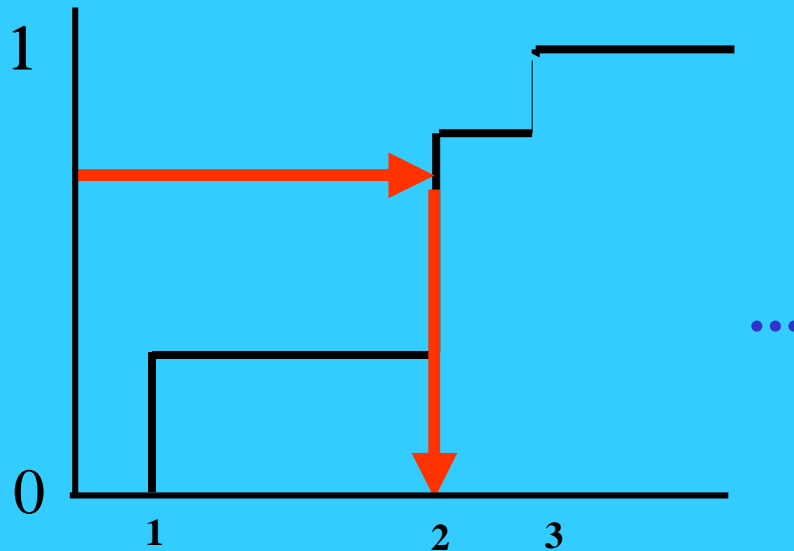
$$p_k^*(\mathbf{u}) = \sum_{\alpha=1}^n \lambda_{\alpha}(\mathbf{u}) \cdot I(u_{\alpha}; k) + [1 - \sum_{\alpha=1}^n \lambda_{\alpha}(\mathbf{u})] \cdot m_k$$

- Determine weights  $\lambda_{\alpha}(\mathbf{u}), \alpha=1, \dots, n$  by the well known “normal system” or kriging.
- Kriging weights account for two things:
  - clustering of the data locations
  - closeness of the data to the location being considered



STEP 3 - (d)

# Draw a Simulated Value



- probabilities  $p_k^*(u)$ ,  $k=1,\dots,K$  are given by kriging
- Procedure:
  - draw a random number  $\in [0,1]$
  - find which class  $k$  is specified by the random number
  - assign  $k$  to node
- Since the conditional probabilities were estimated by kriging with a given variogram  $\gamma_k(\mathbf{h})$ ,  $k=1,\dots,K$ , the simulated values, taken all together, will reproduce those variograms  $\gamma_k(\mathbf{h})$ ,  $k=1,\dots,K$



# Detailed Steps in SISIM

1. Establish grid network and coordinate system ( $Z_{rel}$ -space)
2. Assign data to the nearest grid node (take the closest of multiple data assigned to same node)
3. Determine a random path through all of the grid nodes
  - (a) find nearby data and previously simulated grid nodes
  - (b) construct the conditional probabilities by kriging
  - (c) draw simulated value from conditional distribution
4. Check results
  - (a) honor data?
  - (b) honor global proportions?
  - (c) honor variogram?
  - (d) look reasonable



START OF PARAMETERS:

```

1
5
0.5 1.0 2.5 5.0 10.0
0.12 0.29 0.50 0.74 0.88
../data/cluster.dat
1 2 0 3
direct.ik
1 2 0 3 4 5 6 7
0
0.61 0.54 0.56 0.53 0.29
-1.0e21 1.0e21
0.0 30.0
1 0.0
1 1.0
1 30.0
cluster.dat
3 0
0
sisim.dbg
sisim.out

```

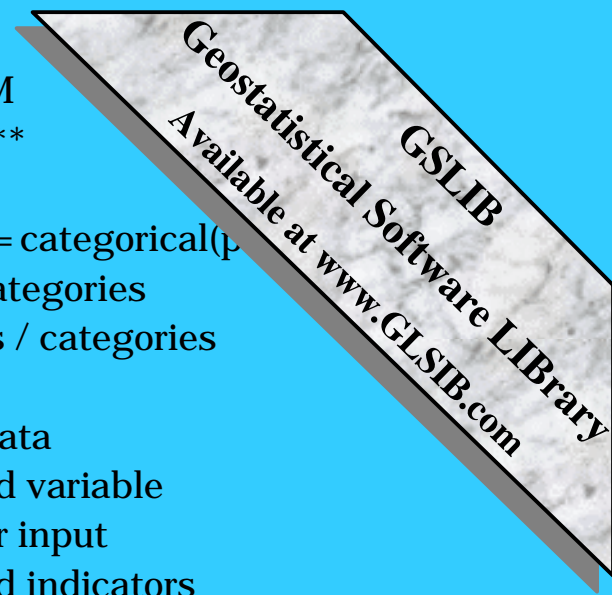
Parameters for SISIM

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```

\ 1= continuous(cdf), 0= categorical(pdf)
\ number thresholds/categories
      \ thresholds / categories
\ global cdf / pdf
      \ file with data
\ columns for X,Y,Z, and variable
\ file with soft indicator input
\ columns for X,Y,Z, and indicators
\ Markov-Bayes simulation (0= no,1= yes)
\ calibration B(z) values
      \ trimming limits
\ minimum and maximum data value
\ lower tail option and parameter
\ middle option and parameter
\ upper tail option and parameter
\ file with tabulated values
\ columns for variable, weight
\ debugging level: 0,1,2,3
\ file for debugging output
\ file for simulation output

```





```
1
50 0.5 1.0
50 0.5 1.0
1 1.0 10.0
69069
12
12
1
0
0 3
0
20.0 20.0 20.0
0.0 0.0 0.0
0 2.5
0
1 0.15
1 0.85 0.0 0.0 0.0
    10.0 10.0 10.0
1 0.10
1 0.90 0.0 0.0 0.0
    10.0 10.0 10.0
1 0.10
1 0.90 0.0 0.0 0.0
    10.0 10.0 10.0
```

```
\ number of realizations
\ nx,xmn,xsiz
\ ny,ymn,ysiz
\ nz,zmn,zsiz
\ random number seed
\ maximum original data for each kriging
\ maximum previous nodes for each kriging
\ maximum soft indicator nodes for kriging
\ assign data to nodes? (0= no,1= yes)
\ multiple grid search? (0= no,1= yes),num
\ maximum per octant (0= not used)
    \ maximum search radii
\ angles for search ellipsoid
\ 0= full IK, 1= median approx. (cutoff)
\ 0= SK, 1= OK
\ One nst, nugget effect
    \ it,cc,ang1,ang2,ang3
    \ a_hmax, a_hmin, a_vert
\ Two nst, nugget effect
    \ it,cc,ang1,ang2,ang3
    \ a_hmax, a_hmin, a_vert
\ Three nst, nugget effect
    \ it,cc,ang1,ang2,ang3
    \ a_hmax, a_hmin, a_vert
```

