# Day 2 Lab Part III

ISI Short Course | June 1-3

### create a 1D SAR weight matrix

Use a handy LatticeKrig function to do it. Note diagonal value muxst be greater than 2.

```
B<- LKDiag( c( -1, 3,-1), 10, full=TRUE)
print( B)</pre>
```

```
##
           [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9] [,10]
##
     [1,]
              3
                   -1
                          0
                                 0
                                       0
                                             0
##
    [2,]
             -1
                    3
                         -1
                                             0
                                                   0
                                                         0
                                                               0
                                                                       0
                                 0
                                       0
##
    [3,]
              0
                   -1
                               -1
                                                                       0
    [4,]
              0
                    0
                                             0
                                                   0
                                                         0
                                                               0
                                                                       0
##
                         -1
                                3
                                      -1
##
    [5,]
              0
                    0
                          0
                                       3
                                            -1
                                                   0
                                                               0
                                                                       0
                               -1
    [6,]
##
              0
                    0
                          0
                                0
                                      -1
                                             3
                                                  -1
                                                         0
                                                               0
                                                                       0
    [7,]
                                                   3
                                                                       0
##
              0
                    0
                          0
                                 0
                                       0
                                            -1
                                                        -1
                                                               0
##
     [8,]
              0
                    0
                          0
                                 0
                                       0
                                             0
                                                  -1
                                                         3
                                                                       0
                                                              -1
##
    [9,]
              0
                    0
                           0
                                 0
                                       0
                                             0
                                                   0
                                                               3
                                                                      -1
                                                        -1
                                                                       3
## [10,]
              0
                    0
                           0
                                 0
                                             0
                                                   0
                                                         0
                                                              -1
                                       0
```

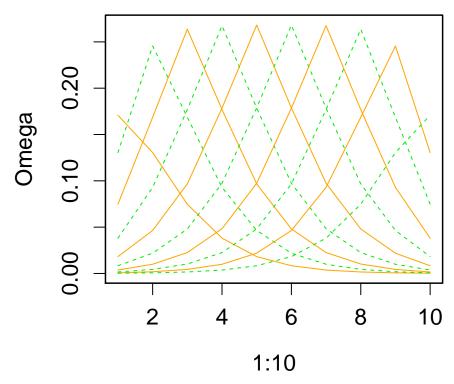
Now find precision matrix and the covariance matrix. The shapes are close to double exponential.

```
fields.style()
Q<- t( B) %*% B
print( Q)</pre>
```

```
[,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9]
##
           [,1]
##
     [1,]
             10
                   -6
                           1
                                 0
                                       0
                                              0
##
    [2,]
             -6
                          -6
                                              0
                                                    0
                                                          0
                                                                0
                                                                        0
                   11
                                 1
                                       0
##
    [3,]
              1
                   -6
                          11
                                -6
                                       1
                                                                        0
    [4,]
##
              0
                     1
                          -6
                                              1
                                                    0
                                                          0
                                                                0
                                                                        0
                                11
                                      -6
##
    [5,]
              0
                     0
                           1
                                -6
                                      11
                                            -6
                                                    1
                                                          0
                                                                0
                                                                        0
##
    [6,]
              0
                     0
                           0
                                 1
                                      -6
                                            11
                                                   -6
                                                          1
                                                                0
                                                                        0
##
    [7,]
              0
                     0
                           0
                                 0
                                       1
                                            -6
                                                   11
                                                         -6
                                                                1
                                                                        0
     [8,]
              0
                     0
                           0
                                       0
                                              1
                                                   -6
##
                                 0
                                                         11
                                                               -6
                                                                        1
##
    [9,]
              0
                     0
                           0
                                 0
                                       0
                                              0
                                                    1
                                                         -6
                                                               11
                                                                       -6
## [10,]
                                 0
                                       0
                                              0
                                                    0
                                                          1
                                                               -6
                                                                       10
```

```
Omega<- solve( Q)
matplot( 1:10, Omega, type="1", col=1:2, lty=1:2)
title("Slices of the covariance matrix")</pre>
```

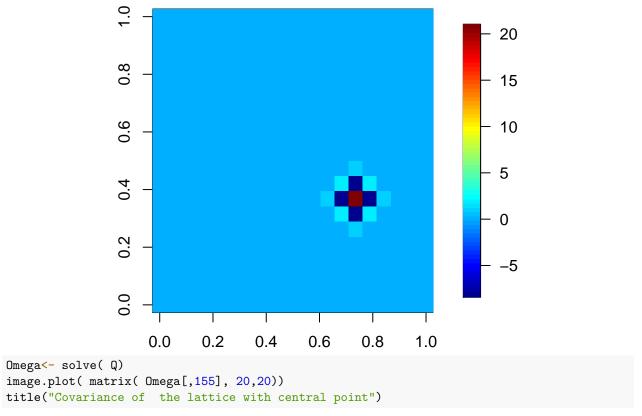
## Slices of the covariance matrix



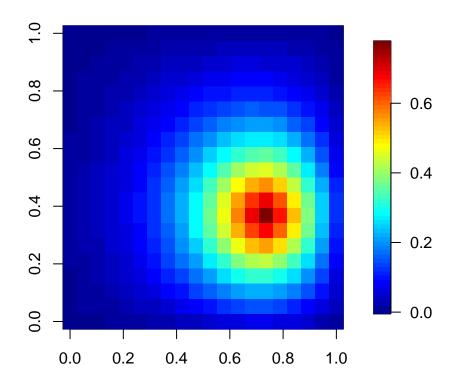
Now use LatticeKrig utilities to look at some MRFs in 2D. This is for a 20X20 lattice. **NC.buffer=0** below means no extra grid points are added at the edges.

```
sDomain<- cbind( c(0,1), c( 0,1))
LKinfo<- LKrigSetup( sDomain, NC= 20, NC.buffer=0, nlevel=1, a.wght=4.1)
Qsparse<- LKrig.precision(LKinfo)
Q<- spam2full( Qsparse)
# check
dim( Q)

## [1] 400 400
# pull off row 155 and reshape to an image.
image.plot( matrix( Q[,155], 20,20))</pre>
```

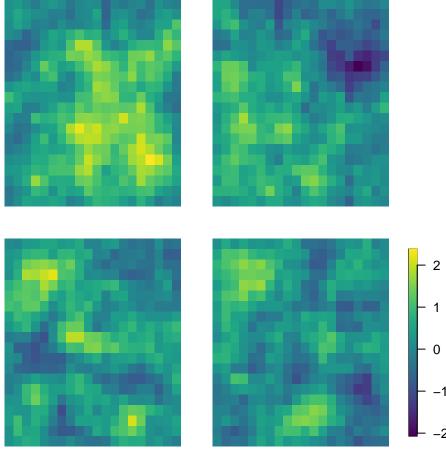


## Covariance of the lattice with central point



#### Simulate some MRFs

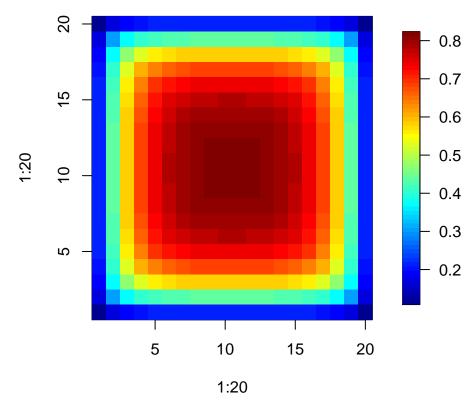
How about 4? Note that the coefficient fields are big vectors. Have to reshape to look at them.



#### Exercises

- 1. How does the covariance at the 155 location change when a.wght=4.01? How about a.wght=10?
- 2. Take a look at the process variances (diag(Omega))

```
image.plot( 1:20, 1:20, matrix( diag( Omega),20,20))
```



You can see some edge effects.

How many rows and columns of the lattice should you ignore so that the variances are approximately constant? Say within about 90%? The LatticeKrig default is 5 on each side.