Some SAR Processes

Short course Large spatial data.

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

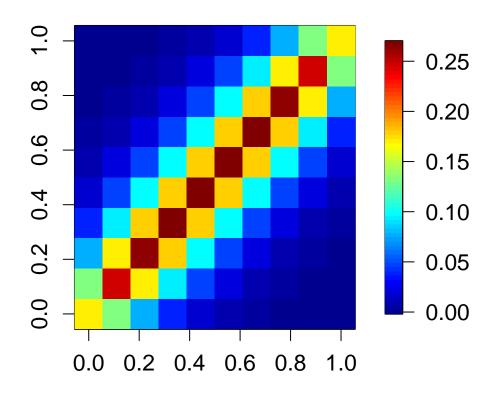
Now find precision matrix and the covariance matrix. The shapes are close to double exponential.

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

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

```
Omega<- solve( Q)
imagePlot( Omega)
title("The covariance matrix")</pre>
```

The covariance matrix



A 2D example

Now use LatticeKrig utilities to look at some SARs in 2D. This is for a 30X30 lattice. **NC.buffer=0** below means no extra grid points are added at the edges. This is not recommended as there are edge effects!

First define the mode with a setup function.

```
sDomain <- cbind( c(0,1), c( 0,1))

LKinfo <- LKrigSetup( sDomain, NC= 30, NC.buffer=0, nlevel=1, a.wght=4.05)

# print(LKinfo) for a summary of the model with its defautl choices.
```

Now use this specification to create some of the model components.

```
Qsparse<- LKrig.precision(LKinfo)
# covert the sparse format to a full one for plotting
Q<- spam2full( Qsparse)
# check
dim( Q)</pre>
```

[1] 900 900

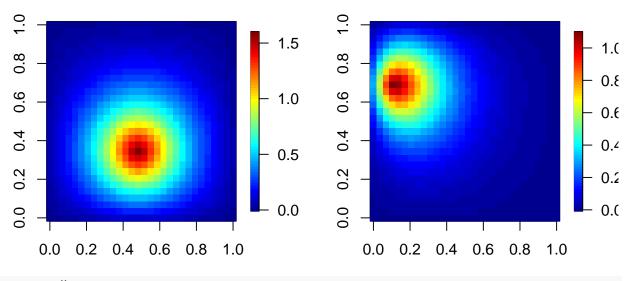
```
Omega<- solve( Q)
set.panel(1,2)</pre>
```

plot window will lay out plots in a 1 by 2 matrix

```
par( pty="s")
#15 + 10*90
# pull off (15,10) and reshape to an image.
image.plot( matrix( Omega[,15+10*30], 30,30))
title("Covariance of the lattice at [15,11]")
image.plot( matrix( Omega[,4+20*30], 30,30))
title("Covariance of the lattice at [4,21]")
```

Covariance of the lattice at [15,11]

Covariance of the lattice at [4,21]



set.panel()

plot window will lay out plots in a 1 by 1 matrix

Simulate some MRFs

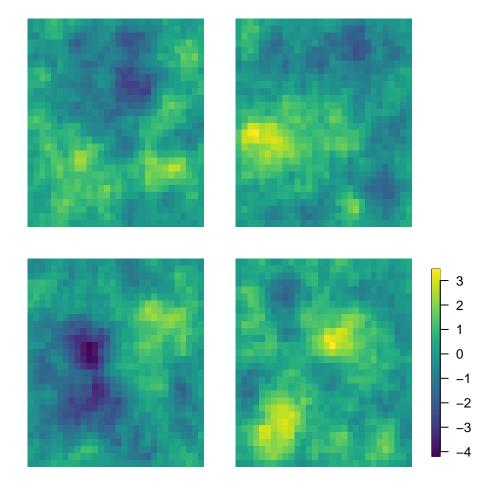
How about 4? Note that the coefficient fields are big vectors. Have to reshape to look at them.

```
set.seed(123)
look <- LKrig.sim( LKinfo=LKinfo, M=4 ,just.coefficients=TRUE )</pre>
```

Now plot them – this takes way more code than generating them!

```
# take a look
set.panel( 2,2)
```

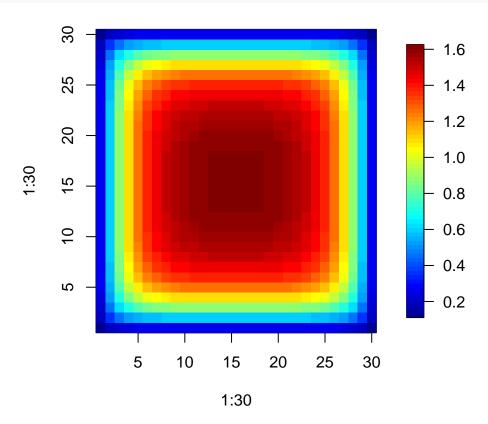
plot window will lay out plots in a 2 by 2 matrix



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(1:30, 1:30, matrix(diag(Omega),30,30))



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.