Chapter 2: Introduction to Slepian Functions

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After having worked through Chapter 1: Introduction to Spherical Harmonics, we will now have a look at what spherical Slepian functions are. The goal of this chapter is not to give a solid foundation of the mathematical intricacies. This can be obtained from the many research articles in the literature. Rather, we will look at some basic ideas and plot the first few Slepian functions.

1 Basic ideas

When plotting spherical harmonics in chapter 1, it became clear that all of these spherical harmonics are functions that cover the entire sphere. This is of course an advantage if we try to describe functions (or maps), that cover the entire planet, such as gravity fields, magnetic fields, etc. We saw in chapter 1 that we can generate pretty much any spatial pattern we like by simply summing up different spherical harmonics each multiplied by a different factor called "coefficient".

But what can we do if we only have information within a specific region? Sometimes it's fine to just describe that information as point values. But in some cases, as for example when we measure gravity or magnetic data at satellite altitude, and we want to calculate, what the corresponding magnetic or gravity field looks like on the planet's surface (assuming there are no gravity or magnetic sources between the planet and the satellite, or that they are negligible), we need to describe these fields in spherical harmonics.

This is where the Slepian functions come in. Slepian functions are linear combinations of spherical harmonics which means that they are constructed by multiplying each spherical harmonic function with a factor (called Slepian coefficient) and then adding them all up. The name Slepian function stems from the first author of a research article, that described the original idea for 1-dimensional functions.

There are different types of Slepian functions that are constructed in different ways but the basic idea is always to solve an optimization problem, in which we try to balance the number of spherical-harmonic functions we use (the maximum spherical-harmonic degree) in the construction of the Slepian functions, and how much they are concentrated within the region that we are interested in.

2 Classical Scalar Slepian Functions

Let's first look at the classical scalar Slepian functions, which are spectrally limited and spatially optimized. These functions are suitable to use when you are limited to information within a specific region as mentioned above, but only when that information is constrained to the surface of the sphere as point values; classical vector Slepian functions will be addressed in a later chapter. The functions will be bandlimited by the maximum spherical harmonic degree L for region of concentration R. We will choose our region to be the continent **Africa** for our first example. The following examples will include a region of an axisymmetric polar cap, a polar ring, and both polar cap and ring again when rotated away from the north pole.

2.1 Calculating Coefficients for a Region

In the Slepian folder start MATLAB and run:

```
demos_chapter_two(2.1)
```

You may also find it helpful to open demos_chapter_two.m in a text editor and browse the script as you read through this tutorial. The demo file is divded into cases for each subsection of this chapter.

The function glmalpha is used to calculate unit-normalized spherical harmonic coefficients given the following required inputs:

```
TH - a desired region (in this example: 'africa')
```

L - the bandwidth, or maximum spherical harmonic degree.

If you find the program running slowly or freezing up in MATLAB due to your machine's computing power, you may need to edit the demo file and set L in glmalpha() to a smaller value, say 5 for example.

Thus the output of glmalpha we will need for this example is:

G - a unitary matrix of localization coefficients.

There are several other inputs and outputs that may and should be used later; see help glmalpha for more details.

We specify one other input to indicate the ordering of the spherical harmonic coefficients. In our codes we use two different methods of ordering: ADDMOUT and ADDMON.

Remember that spherical harmonics have a degree l and an order m. Degrees are all integers greater than zero and orders for each degree vary between -l and l.

We could order these spherical harmonics as

This is called the ADDMOUT ordering and is the ordering in which the columns of the Slepian matrices in glmalpha, gradvecglmalpha, and all these other similar functions are returned.

The other ordering that we may use is:

This one we call ADDMON and it is the ordering in which the coefficients from functions like LocalInnerField (which will be discussed in a later chapter) are returned. Every function that returns coefficients or matrices of coefficients should state in their help menu in which format the coefficients are. Just run help [function] in MATLAB.

Returning to our example, if you run help glmalpha you will read that it uses the ADDMOUT ordering and thus requires us to use 0 as the fourth input. We will choose a value of L=20. Now we can run:

and observe the output G which should be a $(L+1)^2 \times (L+1)^2$ orthonormal matrix of coefficients. Each column of G contains a set of coefficients for a single slepian function which are in the ADDMOUT ordering:

$$G = \begin{bmatrix} g_{00,1} & g_{00,2} & \cdots & g_{00,(L+1)^2} \\ g_{1-1,1} & g_{1-1,2} & \cdots & g_{1-1,(L+1)^2} \\ \vdots & \vdots & & \vdots \\ g_{LL,1} & g_{LL,2} & \cdots & g_{LL,(L+1)2} \end{bmatrix}$$

where the subscript of each element contains the lm values and the slepian function number separated by a comma respectively. The first column G(:,1) is the coefficients corresponding to the "best" Slepian function, or the one that is most spatially concentrated within our chosen region of Africa. G(:,2) would be the second best, G(:,3) would be the third best, and so on until $G(:,(L+1)^2)$.

Once the coefficients have been calculated we desire a neutral type of sorting in the format called LMCOSI to prepare for plotting.

In the lmcosi format, we explicitly write out the degrees, orders, and coefficients which in the following I will call c_{lm}

l	\mathbf{m}	co	si
0	0	c_{00}	0
1	0	c_{10}	0
1	1	c_{1-1}	c_{11}
2	0	c_{20}	0
2	1	c_{2-1}	c_{21}
2	2	c_{2-2}	c_{22}
:	:	:	:

As you can see this last ordering is an entire matrix instead of just a vector. This is a disadvantage to run calculations but it is less ambiguous than just a vector of coefficients.

Luckily we have programs that take care of these issues. These programs have the creative names coef2lmcosi and lmcosi2coef. You can probably guess from the name what each does.

For our Africa example we will need to feed our coefficients into coef2lmcosi and again tell it whether they are in the ADDMOUT or ADDMON format. This is indicated by using the value 1 for the second input (see help coef2lmcosi). The output is saved as *lmcs* but you can name it whatever you would like. Run:

```
lmcs=coef2lmcosi(G(:,1),1)
```

The first column of G is used for now in our example for simplicity's sake.

We now have one more step before we are ready to plot and that is to evaluate the coefficients using the function plm2xyz. The required inputs are our *lmcosi* format matrix and a chosen resolution. You can choose res=1, or 0.5, or even 0.1 for a very high resolution image, but keep in mind this will also use much calculation time and memory. We will give the output the arbitrary name 'data' and choose a resolution of 0.5. Run:

```
data=plm2xyz(lmcs,0.5)
```

And finally to plot we will use plotplm:

```
plotplm(data,[],[],1,0.5)
```

where the fourth input chooses the plotting method. The value '1' corresponds to a Mollweide projection of Earth.

We use kelicol(1) for a more clear color scheme where white sections are of value 0. It is recommended to change the maximum and minimum of the color scheme which you can do by running:

```
caxis([-1,1]*max(abs(caxis)));
```

As expected, the fluctuations are concentrated on the center of Africa and negligible elsewhere. The plot produced by demos_chapter_two(2.1) should match Figure 1.

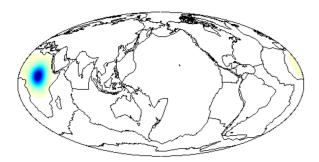


Figure 1: Result of example using glmalpha() with L=20 to calculate coefficients, plm2xyz() to evaluate them, and plotplm() to plot with method 1 - the Mollweide projection. Only the first "best" concentrated Slepian function is plotted. The color scheme is kelicol(1) and max/min values are changed to fit the max/min of data plotted.

Exercise: Pick a different a region, or subtract a region from Africa to use as the domain and plot again. You can edit demos_chapter_two.m or write your own .m file.

2.2 Calculating Coefficients for an Axisymmetric Polar Cap

We now look at the case of an axisymmetric polar cap.

You can run:

```
demos_chapter_two(2.1)
```

which will plot both a Mollweide projection where the region is North polar cap, as well as the same region looking down at the North Pole.

The functions used for this type of region follow the same format as in section 2.1, however we have changed the inputs of glmalpha(). TH is now an opening angle in degrees, which have chosen to be 40, and a value of 1 is passed as the third input to indicate we are creating coefficients for a polar cap of radius TH.

2.3 Calculating Coefficients for a Polar Ring

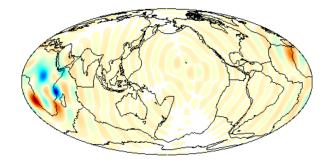
2.4 Rotating a Polar Cap or Ring

3 Linear Combinations of Classical Scalar Slepian Functions

Let's do another quick example where we plot a linear combination of Slepian functions. We can add any combination the columns of G together and also multiply by any value or a vector of scalars. We can repeat the steps from earlier and run:

```
lmcs=coef2lmcosi((5*G(:,1))+(3*G(:,30)),1)
data=plm2xyz(lmcs,0.5)
```

And if you plot it you should see:



Exercsie: Practice plotting different linear combinations of slepian functions. Try multiplying by a vector of random coefficients.

4 Choosing The Number of Slepian Functions

We will now take a look at our vector of eigenvalues, V. You can plot the eigenvalues versus their corresponding slepian function j by setting $J=(L+1)^2$ and running the following:

Using the V from Example One we produce Figure 2 from which we can see that our eigenvalues drop off rapidly in value and become zero near j = 50.

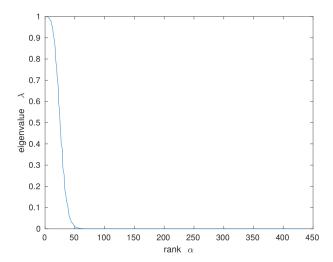


Figure 2: Eigenvalues plotted on the y-axis versus their corresponding Slepian function number on the x-axis. We can see that most of the energy is occurring in functions corresponding to $j \le 50$.

Plotting this curve may be useful in selecting which Slepian functions will best represent the region of interest, as you will soon see.

Lets first plot a linear combination of all solutions and see what it looks like. To evaluate all coefficients you can run:

```
for j=1:J

g\{j\} = plm2xyz(coef2lmcosi(G(:,j),1),.5)

end
```

We can then add the different elements of g into one matrix f:

```
f = zeros*size(g{1}))
for j=1:J
    f = f + g{j}
end
```

and then plot with:

```
plotplm(f,[],[],4,.5)
```

Your figure should match that of Figure 3. We can now see that using all of the eigensolutions does not work when we are trying to define a region like Africa.

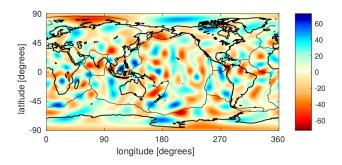


Figure 3: A linear combination of all eigensolutions. L=20.

If we refer back to Figure 2 we can see that our eigenvalues drop off to zero around j = 50. Let's now reset J=50 and see what our plot looks like. You should be able to produce Figure 4, in which we see that our information has been concentrated within our region of choice, Africa.

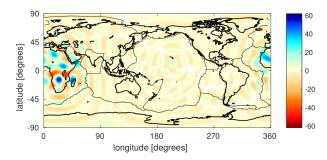


Figure 4: Using a linear combination of the first 50 slepian functions to represent Africa.

Exercise: Create a vector of random numbers the same length as the matrix f we plotted in Figure 4 and multiply it with g{j} inside the for loop. How does this change the plot when using the same eigensolutions? You should produce something similar to Figure 5.

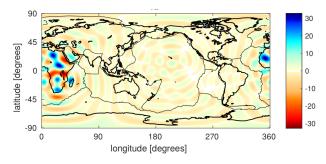


Figure 5: Using a linear combination of the first 50 slepian functions multiplied by a vector of random coefficients.

This tutorial is currently under construction. Please check back later for more by keeping your software updated.