

# User Manual for the paper titled ‘A Wavelet Differentiation Matrix Suite’

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1.

`cascade.m`

The function `cascade.m` computes the scaling function  $\phi$  and the wavelet function  $\psi$  at dyadic rationals. The calling command for this function is

```
>>[x, phi, psi]=cascade(D,q);
```

where `phi` and `psi` stands for  $\phi$  and  $\psi$  respectively.  $D$  and  $q$  are the wavelet genus and the desired dyadic resolution respectively. The implementation goes like this: An inbuilt function in Matlab, `wfilters.m`, computes low pass filter coefficients  $h_k$  and high pass filter coefficients  $g_k$  as

```
>>[h_k, g_k] = wfilters(['db' num2str(D/2)], 'r');
```

Using these  $h_k$ , the matrices  $A_0$  and  $A_1$  are constructed such that

$$\Phi(0) = A_0\Phi(0) \text{ and } \Phi\left(\frac{1}{2}\right) = A_1\Phi(0),$$

where

$$\Phi(x) = [\phi(x)\phi(x+1)\cdots\phi(x+D-1)]^T$$

These matrices are initialized as zero matrices of order  $(D-1)$  using the following Matlab command

```
>>A_0= A_1= zeros(D-1,D-1);
```

Next, the following loop of `cascade.m` constructs the matrices  $A_0$  and  $A_1$

---

```
>>E = hk(1:2:D-1);      %Even numbered low pass filter coefficients h0,h2,h4,...  
>>O = hk(2:2:D);        %Odd numbered low pass filter coefficients h1,h3,h5,...  
>>Firstrow = 1;         % Matrix A0  
>>Lastrow = D/2;
```

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```

>>A0(Firstrow>Lastrow,1)= E;
>>for j=2:2:D-1
    >>Firstrow = Firstrow+1;
    >>Lastrow = Lastrow+1;
    >>A0(Firstrow>Lastrow,j) = 0;
    >>A0(Firstrow>Lastrow,j+1) = E;
>>end;
>>A0 = sqrt(2)*A0;
>>Firstrow = 1;          % Matrix A1
>>Lastrow = D/2;
>>for j=1:2:D-3
    >>A1(Firstrow>Lastrow,j) = 0;
    >>A1(Firstrow>Lastrow,j+1) = E;
    >>Firstrow = Firstrow+1;
    >>Lastrow = Lastrow+1;
>>end;
>>A1(Firstrow>Lastrow,D-1) = 0;
>>A1 = sqrt(2)*A1;

```

---

Now we need to find out the eigenvector of  $A_0$  corresponding to the eigenvalue 1 which is its maximum eigenvalue and this eigenvector is nothing but  $\Phi(0)$ . This is done by following code

```

>>[V,Dia]=eig(A0);          % Get eigenvectors and values of matrix A0
>>[maxeig,index]=max(diag(Dia)); % Find largest eigenvalue (=1)
>>v = V(:,index);          % Pick corresponding eigen vector
>>v = v/sum(v);            % Normalize s.t. sum(v) = 1, (sum(\phi_{k})=1)
>>for k=0:D-2
    >>phi(k*2^q + 1) = v(k+1); % Initialize phi at integers
>>end

```

---

The inbuilt function `eig.m` is used to obtain the eigenvalues and eigenvectors of  $A_0$ . After calculation of  $\Phi(0)$ , `cascade.m` calculate  $\Phi(\frac{1}{2})$ . Last step of `cascade.m` is calculation of  $\psi$  using  $\phi$  using following commands

```

>>psi = zeros(size(phi));
>>if q > 0          % q is required resolution
    >>L2 = L/2;      % L=(D-1)*2^q
    >>phi2 = phi(1:2:L); % Phi at even numerators
    >>first = 1;
    >>for k=1:D
        >>psi(first:first+L2-1) = psi(first:first+L2-1) + sqrt(2)*hp(k)*phi2;
        >>first = first + 2^q/2;
    >>end;
>>else          % q = 0 (a special case)
    >>for m=0:D-1

```

```

>>for k=0:D-1
    >>index = 2*m-k;
    >>if (index >= 0) & (index <= D-1)
        >>psi(m+1) = psi(m+1) + sqrt(2)*hp(k+1)*phi(index+1);
    >>end;
>>end;
>>end;
>>end;

```

---

2.

`dstmat.m`, `dst.m` and `idst.m`

- The `dstmat.m` computes the matrix  $T_{r,j}$  required for discrete scaling function transformation and inverse discrete scaling function transformation.
- The `dst.m` and `idst.m` perform the required calculations for discrete and inverse discrete scaling function transformation respectively.

The function `dstmat.m` first initializes the required matrix  $T_{r,j}$  using

```
>> T = spalloc(K*2^r,K*2^j,(D-1)*2^(r-j));
```

The `spalloc.m` is an inbuilt Matlab routine, `spalloc(M,N,Nz)` creates an  $M \times N$  sparse matrix with all entries zero and with room to eventually hold  $Nz$  non zero elements. After initialization step, following is the main loop of `dstmat.m`

---

```

>>N=K*2^r;                % Number of samples
>>P=K*2^j;                % Number of coefficients
>>Q=2^q;                  % The resolution of phi
>>S = P*Q/N;              % step between values needed in phi is 2^(j-r+q)
>>phi_s = phi(1:S:length(phi));
>>len    = length(phi_s);
>>if len < N                % phi_s fits into a column
    >>for l=0:P-1
        >>firstrow = l*N/P;
        >>lastrow  = firstrow + len-1;
        >>wrap     = lastrow - (N-1);
        >>if wrap > 0,
            >>lastrow = lastrow - wrap;
            >>T(0+shift:wrap-1+shift, l+shift) = phi_s(len-wrap+1:len);
        >>else
            >>wrap=0;
        >>end;
        >>T(firstrow+shift:lastrow+shift, l+shift) = phi_s(1:len-wrap);
    >>end;
>>else
    >>for l=0:P-1            % For each column
        >>firstrow = rem(l*N/P,N);

```

```

>>for k=0:len-1          % For element in phi_s
>>row = rem(firstrow + k, N);
>>T(row+shift, l+shift) = T(row+shift, l+shift) + phi_s(k+shift);
>>end;
>>end;
>>end;
>>T=2^(j/2)*T;

```

---

Both `dst.m` and `idst.m` call `dstmat.m` using the command

```
>> T=dstmat(wavelet,r,j,q,K);
```

The calling commands for `dst.m` and `idst.m` are

```

>>c=dst(f, D);
>>f=idst(c, D, q);

```

Main commands for `dst.m` are:

```

>>T = dstmat(D,r,j,q,L);
>>c = T\f;

```

---

3.

`moments.m` and `tmoments.m`

- The function `moments.m` is called by the command

```
>>moments = moments (h_k, pmax);
```

where `moments` and `h_k` stands for  $M_0^p$  and vector containing the low pass filter coefficients respectively, `pmax` is the number of moments to be computed.

- The function `tmoments.m` computes the translated moments  $M_l^p$ .

4.

`conn.m`, `gal_difmatrix_periodic.m` and `gal_diff_periodic.m`

- The function `conn.m` implements the algorithm for the computation of the connection coefficients. The calling command for `conn.m` is

```
>>Gamma = conn (d, D);
```

where `Gamma` stands for  $\Gamma^d$ ,  $d$  is the order of differentiation and  $D$  is the wavelet genus. First of all it uses the Matlab inbuilt function `wfilters` to calculate low pass and high pass filter coefficients. Then using these low pass filter coefficients it constructs the matrix `A` of the equation

$$(\mathbf{A} - 2^{-d}\mathbf{I})\Gamma^d = 0. \quad (1)$$

Next it uses the following command

```
>>A = A - eye(M)/2^(d);
```

The `eye` is an inbuilt function to construct the identity matrix. It then uses Matlab inbuilt function `svd` to find the solution of (1). The command

```
[U,S,V] = svd(A);
```

produces a diagonal matrix  $S$ , of the same size as that of  $A$  and with nonnegative diagonal elements in decreasing order, and unitary matrices  $U$  and  $V$  such that  $A = USV'$ . The  $(2D - 3)^{th}$  column of  $V$  will serve the purpose i.e. it is the required solution of (1). The normalization is done by following set of commands

---

```
>>if d == 0
    >>Mrow = ones (1,2D-3);
>>else
    >>Mom = moments (hk, d);
    >>Mrow = zeros (1,2D-3);
    >>for c=1:2D-3,
        >>l = c-(D-1);
        >>TMom = tmoments (Mom,l);
        >>Mrow(c) = TMom(d);
    >>end
>>end;
>>c = (-1)^d * fac(d) / (Mrow * V(:,M));
>>Gamma = c * V(:,M);
```

---

- The function `gal_difmatrix_periodic.m` is the Matlab function in the suite which constructs the differentiation matrix for the Galerkin approach in periodic case. The calling command for it is

```
>>difmatrix=gal_difmatrix_periodic(d,N,L,D);
```

where `difmatrix` stands for  $\mathcal{D}^{(d)}$ ,  $d$  is the order of differentiation,  $N = 2^j$  is the size of differentiation matrix,  $L$  is the period of the function which is to be differentiated and  $D$  is the wavelet genus.

- The function `gal_diff_periodic.m` will differentiate a given function  $f$  using the differentiation matrix produced by `gal_difmatrix_periodic.m`. The calling command is

```
>>[derivative, error] = gal_diff_periodic(j,d,D,L);
```

where `derivative` and `error` stands for  $\mathcal{D}^{(d)}f$  and  $E^{(d)}(f,j)$  respectively.  $L$  is the period of the function to be differentiated. This function will ask for the function to be differentiated.

5.

`cascade_der.m`, `collo_difmatrix_periodic.m` and `collo_diff_periodic.m`

- The function `cascade_der.m` computes the values of  $\phi^{(d)}$  and  $\psi^{(d)}$  at dyadic rationals. The calling command for it is

```
>>[x, phi_d, psi_d]=cascade_der(D,q,d);
```

where `phi_d` and `psi_d` stands for  $\phi^{(d)}(x)$  and  $\psi^{(d)}(x)$  respectively. The `moments.m` and `tmoments.m` are called by `cascade_der.m`.

- Matlab function `collo_difmatrix_periodic.m` calls `cascade_der.m` and then constructs the matrix  $\mathcal{D}^{(d)}$  in case of collocation approach in periodic case. The calling command for `collo_difmatrix_periodic.m` is

```
T=collo_difmatrix_periodic(D,r,j,q,d);
```

- The function `collo_diff_periodic.m` will differentiate a given function  $f$  using the differentiation matrix produced by `collo_difmatrix_periodic.m`. The calling command for it is

```
>>[derivative, error] = collo_diff_periodic(j,D,L,d);
```

where `derivative` and `error` stands for  $\mathcal{D}^{(d)}$  and  $E^{(d)}(f, j)$  respectively. It will ask for a function to be differentiated and will return the value of derivative of the function at the nodal points and the error.

## 6.

`L_daubfilt.m` and `R_daubfilt.m`

- The function `L_daubfilt.m` computes the left low pass filter coefficients  $h^L$ . The calling command is:

```
>>[h_L]=L_daubfilt(k,D);
```

where `h_L` stands for  $h^L$ .

- The function `R_daubfilt.m` computes the right low pass filter coefficients  $h^R$ . The calling command is:

```
>>[h_R]=R_daubfilt(k,D);
```

where `h_R` stands for  $h^R$ .

## 7.

`L_moments.m` and `R_moments.m`

- The function `L_moments.m` calculates the moments of left hand side boundary scaling functions. The calling command is

```
>>[moments_Lp]=L_moments(D,p);
```

where `moments_Lp` stands for  $m_k^{L,p} = \{m_k^{L,p}, \quad k = 0, \dots, M-1\}$ .

- The function `R_moments.m` calculates the moments of right hand side boundary scaling functions. The calling command for it is

```
>>[moments_Rp]=R_moments(D,p);
```

8.

`dstmat_nonper.m`

The function `dstmat_nonper.m` constructs the quadrature matrix  $C$  in non-periodic case of Galerkin approach. The calling command is

```
>>[C]=dstmat_nonper(D,N);
```

The `dstmat_nonper.m` calls the functions `L_momoments.m`, `R_moments.m`. The main Matlab commands used in `dstmat_nonper.m` are

---

```
>>ind=0;
>>st=0;
>>for k=0:N-1
    >>if(k>=0 & k<=M-1)
        >>for m=0:M-1
            >>for l=0:M-1
                >>a(m+1,l+1)=x_1(k+l+1)^(m);
            >>end
            >>[L_mo]=L_moments(D,k);
            >>b(m+1)=L_mo(m+1);
        >>end
        >>sol=a\b' ;
        >>for j=0:M-1
            >>C(k+1,j+1)=sol(j+1);
        >>end
    >>elseif((k>=M & k<=N-2*M+1))
        >>for m=0:M-1
            >>for l=0:M-1
                >>a(m+1,l+1)=l^m;
            >>end
            >>[L_mo]=mom1(D);
            >>b(m+1)=L_mo(m+1);
        >>end
        >>sol=a\b' ;
        >>for j=M+ind:M+ind+(M-1)
            >>C(k+1,j+1)=sol(j-(M+ind)+1);
        >>end
        >>ind=ind+1;
    >>else
        >>for m=0:M-1
            >>for l=0:M-1
                >>a(m+1,l+1)=x_1(N-M+l+1)^(m);
            >>end
            >>[R_mo]=R_mom(D,st);
            >>b(m+1)=R_mo(m+1);
        >>end
        >>if(st<M-1)
            >>st=st+1;
```

```

>>end
>>sol=a\b' ;
>>for j=N-M:N-1
>>C(k+1,j+1)=sol(j-(N-M)+1);
>>end
>>end
>>end

```

---

9.

L\_alpha.m and R\_alpha.m

- The function L\_alpha.m implements the algorithm for the calculation of  $\alpha_{m,i}^L$ . The calling command is

```
>> [alpha_L_mi]=L_alpha(m,i,D,N);
```

where alpha\_L\_mi stands for  $\alpha_{m,i}^L$ ,  $D$  is wavelet genus and  $N$  is the dimension of the subspace i.e.  $2^j$ . The functions L\_firstsum\_alpha.m and conn.m are called by L\_alpha.m.

- The Matlab function R\_alpha.m (which calls R\_firstsum\_alpha.m and conn.m) works on the same line as its left counterpart. The calling command is

```
>> [alpha_R_mi]=R_alpha(m,i,D,N);
```

where alpha\_R\_mi stands for  $\alpha_{m,i}^R$ .

10.

L\_phi.m, L\_phi\_origin.m, L\_ro.m, R\_phi.m, R\_phi\_origin.m and R\_ro.m

- The function L\_phi.m computes the value of  $\phi_k^L(x)$  for any  $x$  other than 0. The calling command is

```
>>[phi_Lk]=L_phi(k,x,D,h,supp,phi);
```

where phi\_Lk stands for  $\phi_k^L(x)$ , and h, phi and supp are obtained using the function cascade.m as follows

```

>>[y,phi,psi]=cascade(D,q);
>>h=y(2)-y(1);
>>supp=(D-1)*(1/h)+1;

```

The functions L\_daubfilt.m and L\_firstsum\_phi.m are called by L\_phi.m.

- The function L\_phi\_origin.m calculates  $\phi_k^L(0)$  and  $\rho_{k,k}^L$  for  $D = 4$  and  $D = 8$ . The calling command for this function is

```
>>[phi_Lk0, rho_Lkk]=L_phi_origin(k,D,h,supp,phi);
```

where phi\_Lk0 and rho\_Lkk stands for  $\phi_k^L(0)$  and  $\rho_{k,k}^L$  respectively. The function L\_phi\_origin.m for  $D = 4$  uses following Matlab commands



```

>>x(1)=1/2;
>>x(2)=1;
>>y(1)=L_phi(k,x(1),D,h,supp,phi);
>>y(2)=L_phi(k,x(2),D,h,supp,phi);
>>p=polyfit(x,y,N-1);
>>x1=0;
>>y1=polyval(p,x1);
>>ro=-(y1^2)/2;

```

It is clear from above that `L_phi_origin.m` calls the function `L_phi.m` two times to calculate  $\phi_k^L(1/2)$  and  $\phi_k^L(1)$ . Moreover, it uses the inbuilt Matlab function `polyfit` to interpolate the function  $\phi_k^L(x)$  and then use the inbuilt function `polyval` to calculate  $\phi_k^L(0)$ . The Matlab function `polyfit` is called by the command

```
>>P = polyfit(X,Y,N);
```

It finds the coefficients of a polynomial  $P(X)$  of degree  $N$  that fits the data  $Y$  best in a least-squares sense.  $P$  is a row vector of length  $N+1$  containing the polynomial coefficients in descending powers. The Matlab function `polyval` is called by the command

```
>>Y = polyval(P,X);
```

It returns the value of a polynomial  $P$  evaluated at  $X$ .  $P$  is a vector of length  $N+1$  whose elements are the coefficients of the polynomial in descending powers.

- For the calculations of  $\rho_{k,p}^L$  when  $k \neq p$ , the function `L_ro.m` is available in the suite. The calling command for it is

```
>>[rho_Lkp]=L_ro(D,L,h,supp,phi);
```

where `rho_Lkp` stands for the vector  $\{\rho_{k,p}^L | k \neq p\}$ . The functions `L_partialsum_ro.m`, `L_daubfilt.m`, `conn.m`, `L_phi_origin.m` and `L_alpha.m` are called by `L_ro.m`.

- The Matlab functions `R_phi.m`, `R_phi_origin.m` and `R_ro.m` are available in the suite which all work on the similar lines as their left counterparts.

11.

`gal_difmatrix_nonper.m` and `gal_diff_nonper.m`

- The function `gal_difmatrix_nonper.m` constructs the differentiation projection matrix  $D^{(1)}$  in non-periodic case of Galerkin approach. The calling command is

```
>>[difmatrix]=gal_difmatrix_nonper(D,N);
```

where `difmatrix` stands for  $\mathcal{D}^{(1)}$ . The function `gal_difmatrix_nonper.m` has two for loops for the row index  $l$  and column index  $k$ . The main Matlab commands of the function `gal_difmatrix_nonper.m` are

```

L_sol=L_ro(D,L,h,supp,phi);
R_sol= R_ro(D,L,h,supp,phi);
con_1=conn(1,D);
>>for l=0:N-1
    >>for k=0:N-1
        *****
>>if(l>=0 & l<=M-1)
        *****
>>if(k>=0 & k<=M-1)
    >>if(k==l)
        >>[y,ro]=L_phi2(k,D,h,supp,phi) ;
        >>D(l+1,k+1)=ro;
    >>else
        >>ind=(l)*M+k;
        >>D(l+1,k+1)=L_sol(ind);
    >>end
>>elseif(k>=M & k<=N-2*M+1)
    >>D(l+1,k+1)=L_alpha(k,l,D,L);
>>end
        *****
>>elseif(l>=M & l<=N-2*M+1)
        *****
>>if(k>=0 & k<=M-1)
    >>D(l+1,k+1)=-L_alpha(l,k,D,L);
>>elseif(k>=M & k<=(N-2*M+1))
    >>ind=l-k+(D-1);
    >>if(ind>=1 & ind<=2*(D-2)+1)
        >>D(l+1,k+1)=con_1(ind);
    >>end
>>elseif(k>=(N-2*M+2) & k<=N-1)
    >>if(-(l-(N-2*M+1))>=0 & -(l-(N-2*M+1))<=M-1...
        & -(k-(N-1))>=0 & -(k-(N-1))<=M-1)
        >>D(l+1,k+1)=-R_alpha(-(l-(N-1)),-(k-(N-1)),D,L);
    >>else
        >>D(l+1,k+1)=0;
    >>end
>>end
        *****
>>elseif(l>=(N-2*M+2) & l<=N-1)
        *****
>>if(k>=M & k<=N-2*M+1)
    >>if(-(l-(N-1))>=0 & -(l-(N-1))<=M-1 & -(k-(N-2*M+1))>=0...
        & -(k-(N-2*M+1))<=M-1)
        >>D(l+1,k+1)=R_alpha(-(k-(N-1)),-(l-(N-1)),D,L);
    >>else
        >>D(l+1,k+1)=0;

```

```

>>end
>>elseif(k>=(N-2*M+2) & k<=N-1)
>>if(k==1)
>>[y,ro]=R_phi_origin(-(k-(N-1)),D,h,supp,phi);
>>D(l+1,k+1)=ro;
>>else
>>ind=(1-(N-2*M+2))*(2*M-2)+(k-(N-2*M+2));
>>if(ind<=length(R_sol))
>>D(l+1,k+1)=R_sol(ind);
>>end
>>end
>>end
>>end
*****
>>end
*****
>>end
>>end

```

---

- The function `gal_diff_nonper.m` is included in our suite to construct the differentiation matrix  $\mathcal{D}^{(1)} = C^{-1}D^{(1)}C$  in the non-periodic case of Galerkin approach. The calling command for it is

```
>>[difmatrix]=gal_diff_nonper(D,N);
```

where `difmatrix` stands for  $\mathcal{D}^{(1)}$ . The Matlab commands for `gal_diff_nonper.m` are

---

```

function [difmatrix]=gal_diff_nonper(D,N)
D_matrix=gal_difmatrix_nonper(D,N);
C=dstmat_nonper(D,N);
difmatrix=inv(C)*D_matrix*C;

```

---

12.

`Bn_spline.m`

The calling command for the function `Bn_spline.m` is

```
>>[beta_nxk]=Bn_spline(x,n,k);
```

where `beta_nxk` stands for  $\beta^n(x - k)$ . The code is

---

```

>>function [beta]=Bn_spline(x,n,k)
>>x=x-k;
>>sum=0;
>>for j=0:n+1
>>    if(j~=0)
>>        combos=combnats(1:n+1,j);
>>        tem=size(combos,1);

```

```

        >>else
            >>tem=1;
        >>end
        >>sum=sum+(((−1)^j)/factorial(n))*tem*pow_fun(n,x+((n+1)/2)−j);
    >>end
    >>beta=sum;
    >>beta=beta*factorial(n);

```

---

This function evaluates the value of  $\beta^n(x-k)$ . The function `pow_fun.m` is available in the suite which computes  $[x]_+^n = \max\{0, x\}^n$ .