```
matlab.mixin.indexing.RedefinesBrace
   properties
        Phasor3D %A 3D array reprensenting the harmonics of PhasorArray. Can
be a 3D Array of double, of sim or sdpvar
    end
    methods
        function obj = PhasorArray(varargin, varg)
            %Contructor for PhasorArray
            % Synthax :
                pA = PhasorArray(A) where A is 3D array of
                    double/sym/sdpvar, with size(A,3) is odd.
                pA = PhasorArray(A, "z pospart", true) where A is 3D array
                    of double/sym/sdp, with A(:,:,1) represent the zero
phasor,
                    A(:,:,k) is the k+1 phasor. Negative phasor are deduced
as
            응
                    conjugate of provided phasors.
            응
                pA = PhasorArray(A0, Ap, "z pospart", true) where A0 is a
            응
                    matrix and Ap a 3D Array of compatible size, AO contain
the
                    O phasor, Ap the positive phasor. Negative phasor are
                    deduced by conjugaison.
                pA = PhasorArray(A, "reduce", true) , delete symmetrically
extremal slice
                    in A with value 0.
            %Expect a 3D array of double, sim or sdp var
            % optionnal argument include
               -"reduce" (true/false) to automatically delete extremal 0
            응
                    phasor
                -"z pospart" for zero positive part, in case only 0 and
            응
            응
                    positive harmonics are given, and negative harmonics are
                    deduce by taking conjugate of positive harmonics (for
            응
real
            응
                    valued periodic matrices)
            응
                optionnal acceptable input
                PhasorArray(n,m) is the same as PhasorArray(zeros(n,m))
                PhasorArray(X0, Xpos, "z pospart", true) is an alternate way
                to provide 0 pos part phasor
            arguments (Repeating)
                varargin
            end
            arguments
                varg.reduce=false
                varg.z pospart=false
            end
            if nargin == 0
                obj = PhasorArray(1);
                return
```

classdef PhasorArray < matlab.mixin.indexing.RedefinesParen &</pre>

```
end
            if varg.z pospart %si on a indiqué fournir un 0-phaosr et les
phasor positif
                if numel(varargin)==1
                    varP=varargin{1};
                    varP0=varP(:,:,1);
                    varPpos=varP(:,:,2:end);
                    obj = PosPart2PhasorArray(varP0, varPpos);
                else
                    assert(numel(varargin) == 2, 'If 0 and PosPhasor are
provided, only 2 argument must be provided')
                    obj = PosPart2PhasorArray(varargin{:});
                end
            elseif numel(varargin)>1 %SINON, si on a fourni plusieurs
argument scalaire, c'est qu'on fournit une taille pour un phasor array nul
assert(and(isscalar(varargin{1}),isscalar(varargin{2})),"error")
                obj.Phasor3D=zeros(varargin{:});
                varg.reduce =0;
            elseif isa(varargin{1},'PhasorArray') %si le premier arg est un
phasor array, tous les arguments suivant sont ignorés
                obj.Phasor3D = pvalue(varargin{1});
            elseif isa(varargin{1},'sparsePhasorArray') %si le premier arg
est un sparse phasor array, tous les arguments suivant sont ignorés
                obj = varargin{1}.toPhasorArray();
                return
            else
                assert (mod(size(varargin\{1\},3),2)==1,'Dim3 of 3D array
cannot be even')
                obj.Phasor3D = varargin{1};
            end
            if varg.reduce
                if isa(obj.Phasor3D,'double')
                    obj.Phasor3D=ReduceArray(obj.Phasor3D);
                end
            end
        end
        function info = info(o1, tol)
        % INFO Retrieve and evaluate properties of a PhasorArray object
            info = INFO(o1, tol) computes multiple characteristics of the
            PhasorArray, including harmonic information, real and imaginary
energy
            distribution, and shape properties. It returns a structured
output
        응
            containing these metrics.
        응
        응
            INPUTS:
        응
                   - The PhasorArray object to analyze.
                tol - (Optional) A tolerance used to determine if the object
                      is considered real. Defaults to machine precision if
not
                      specified.
```

```
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            OUTPUT:
                info - A struct containing:
                                         : Logical flag indicating if the
array
                                           is real within the specified
tolerance.
                   - maxImagCoeff
                                         : Maximum magnitude of the imaginary
        응
                                           coefficients.
        응
                   - minImagCoeff
                                         : Minimum magnitude of the imaginary
        응
                                           coefficients.
        응
                   - maxRealCoeff
                                         : Maximum magnitude of the real
                                           coefficients.
        응
                                         : Minimum magnitude of the real
        응
                   - minRealCoeff
                                           coefficients.
                   - iscomplex
                                         : Logical flag indicating if the
array
                                           has non-zero imaginary content.
        응
                                         : Element-wise real energy
                   - realEnergy EW
distribution.
                   imagEnergy EW
                                        : Element-wise imaginary energy
distribution.
        응
                   - realEnergy
                                        : Total real energy.
        응
                   - imagEnergy
                                        : Total imaginary energy.
                   - Energy EW
                                         : Element-wise total energy
distribution.
        용
                   - Energy
                                         : Total energy.
                   - realRelativeEnergy : Fraction of total energy carried
by real
                                           parts.
        응
                   - imagRelativeEnergy : Fraction of total energy carried by
                                           imaginary parts.
                   - imagRelativeEnergy EW : Element-wise relative imaginary
energy.
                   - realRelativeEnergy EW : Element-wise relative real
energy.
        응
                   - h
                                         : Number of harmonics in the array.
        응
                   - size
                                         : Size of the PhasorArray.
                   - isSymmetric
                                         : Logical flag for symmetry check.
        응
                   - isHermitian
                                        : Logical flag for Hermitian check.
                                        : Logical flag for square dimension
                   - isSquare
check.
                                        : Logical flag for scalar dimension
                   - isScalar
check.
                   - isVector
                                         : Logical flag for vector dimension
check.
        응
            NOTE:
                If the array contains small imaginary components but is
flagged
                as real, proceed with caution if the imaginary energy is non-
zero.
            arguments
                01
```

```
tol = []
            end
            if nargin == 1
                warning ("tol is not specified, default to eps, real
evaluation of the PhasorArray may be affected")
            %display information about the PhasorArray
            disp("PhasorArray of size "+num2str(size(o1))+" with
"+num2str(o1.h)+" harmonics")
            [r, \sim] = isreal(o1, tol);
            info.isReal = r;
            info.maxImagCoeff = max(abs(value(mimag(o1))),[],'all');
            info.minImagCoeff = min(abs(value(mimag(o1))),[],'all');
            info.maxRealCoeff = max(abs(value(mreal(o1))),[],'all');
            info.minRealCoeff = min(abs(value(mreal(o1))),[],'all');
            info.iscomplex = \sim r;
            ro1 = mreal(o1);
            info.realEnergy EW = sum(ro1.value.*conj(ro1.value),3);
            info.imagEnergy EW =
sum(value(mimag(o1)).*conj(value(mimag(o1))),3);
            info.imagEnergy = sum(info.imagEnergy EW, 'all');
            info.realEnergy = sum(info.realEnergy EW, 'all');
            info.Energy EW = info.realEnergy EW + info.imagEnergy EW;
            info.Energy = sum(info.Energy EW, 'all');
            info.realRelativeEnergy = info.realEnergy/info.Energy;
            info.imagRelativeEnergy = info.imagEnergy/info.Energy;
            info.imagRelativeEnergy EW = info.imagEnergy EW./info.Energy EW;
            info.realRelativeEnergy EW = info.realEnergy EW./info.Energy EW;
            info.h = o1.h;
            info.size = size(o1);
            info.isSymmetric = issymmetric(o1);
            info.isHermitian = ishermitian(o1);
            info.isSquare = issquare(o1);
            info.isScalar = isscalar(o1);
            info.isVector = isvector(o1);
            if info.imagEnergy>0 && info.isReal
                disp("PhasorArray is real within machine precision, but has
non zero imaginary energy, proceed with caution")
            end
        end
        function [Eew, E] = realEnergy(01)
        %REALENERGY Compute the total real energy of the PhasorArray.
            [Eew,E] = REALENERGY(o1) returns element-wise real energy Eew
            and total real energy E.
            Eew = sum(o1.mreal.value.*conj(o1.mreal.value),3);
            E = sum(Eew, 'all');
        end
```

```
function [Eew, E] = imagEnergy(o1)
            %IMAGENERGY Compute the total imaginary energy of the
PhasorArray.
                [Eew, E] = IMAGENERGY(01) returns element-wise imaginary
energy Eew
            % and total imaginary energy E.
            Eew = sum(o1.mimag.value.*conj(o1.mimag.value),3);
            E = sum(Eew, 'all');
        end
        function [Eew, E] = energy(o1)
            %ENERGY Compute the total energy of the PhasorArray.
               [Eew,E] = ENERGY(o1) returns element-wise energy Eew
            % and total energy E.
            Eew = sum(o1.value.*conj(o1.value),3);
            E = sum(Eew, 'all');
        end
        function [Eew,E] = pageEnergy(o1, normalized, cumulative, plotVar)
            %PAGEENERGY Computes and optionally plots the phasor energy in a
PhasorArray object.
            응
               [Eew, E] = pageEnergy(o1, normalized, cumulative, plotVar)
            응
                This function calculates the energy contributed by each
phasor in the
                PhasorArray object "o1", storing the element-wise energy in
"Eew" and
               the total energy in "E". When the "normalized" flag is set
to true, both
                outputs are normalized by their respective sums. The
"cumulative"
                argument specifies whether the energies should be
accumulated in forward
                ("cumulative") or reverse ("reverse") order, or left
unmodified ("none").
                The "plotVar" argument determines whether to visualize the
results in
               linear scale ("linear"), logarithmic scale ("log"), or skip
plotting
               ("none").
            응
            응
                Inputs:
                                - PhasorArray object containing phasors
                    01
            응
                    normalized - Logical flag to normalize outputs
(default: false)
                    cumulative - String indicating cumulative summation
mode:
                                  'cumulative', 'reverse', or 'none'
(default: 'none')
                                - String for plotting options: 'none',
                    plotVar
'linear', or
```

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                                    'log' (default: 'none')
            응
            응
                Outputs:
            응
                     Eew
                                 - Element-wise phasor energy (matrix)
            응
                     \mathbf{E}
                                 - Total energy of each phasor (vector)
            응
            응
                Example Usage:
                     % Compute and plot phasor energies in linear scale:
            응
                     [Eew, E] = pageEnergy(o1, false, 'none', 'linear');
            arguments
                01
                normalized = false
                cumulative {mustBeMember(cumulative,
{'cumulative', 'reverse', 'none'})} = 'none'
                plotVar {mustBeMember(plotVar, {'none', 'linear', 'log'}) } =
"none"
            end
            for hi = o1.h:-1:0
                oi = o1.extract(hi);
                 [Eew(:,:,hi+1),E(hi+1)] = oi.energy;
            end
            if normalized
                 [EewTotal, Etotal] = o1.energy;
                Eew = Eew./EewTotal;
                E = E./Etotal;
                E(isnan(E))=0;
                Eew(isnan(Eew))=0;
            end
            switch cumulative
                 case "cumulative"
                    Eew = cumsum(Eew, 3);
                     E = cumsum(E);
                case "reverse"
                     Eew = flip(Eew, 3);
                    E = flip(E);
                    Eew = cumsum(Eew, 3);
                    E = cumsum(E);
                     Eew = flip(Eew, 3);
                    E = flip(E);
                case "none"
                otherwise
                     error("cumulative must be 'cumulative', 'reverse' or
'none'")
            end
            if strcmp(plotVar, "linear") || strcmp(plotVar, "log")
                if normalized
                    prefix = "Normalized";
                else
                     prefix = "";
                end
                 switch cumulative
                     case "cumulative"
                    plot(0:o1.h, E(end)-E)
```

```
title("Cumulative " +prefix+"energy of each phasor
(from 0 to harmonic order), 1-Energy")
                    case "reverse"
                    plot(0:o1.h,E)
                        title ("Cumulative " +prefix+"energy of each phasor
in reverse order (from end to harmonic order)")
                    case "none"
                    plot(0:o1.h,E)
                        title(prefix+" Energy of each phasor")
                end
                xlabel("Harmonic order")
                vlabel("Energy")
                if strcmp(plotVar, "log")
                    set(gca, 'YScale', 'log')
                end
            end
        end
        function r = plus(o1,o2)
            % PLUS Overloads the plus (+) operator for the PhasorArray class.
                 PLUS(A,B) returns a new PhasorArray representing the
element-wise
                 sum of A(t) and B(t). The second argument can be a scalar
or a
                 scalar PhasorArray, which is broadcast to match the
dimensions of A.
            if isscalar(o2)
                o2=ones(size(o1,1),size(o1,2))*o2;
            r = PhasorArray(PhasorArrayAdd(o1,o2));
        end
        function r = pageplus(01,02)
            PAGEPLUS(A,B) : A+B = A(t)+B(t)
            %restricted to phasor of same third dimension
            %shouldn't be used directly, use PLUS instead
            %See also PLUS
            r=o1.value+o2.value;
        end
        function r = minus(o1, o2)
            %MINUS overloading for PhasorArray
            %MINUS(A,B) : A-B = A(t)-B(t)
            %MINUS can accept scalar or scalar PhasorArray as second argument
            r = o1 + (-o2);
        end
        function r = uminus(o1)
            %UMINUS overloading for PhasorArray
            %UMINUS(A) : -A = -A(t)
            r = PhasorArray(-pvalue(o1));
        end
```

```
function r = uplus(01)
            %operator overloading for consistency, does nothing
            r = o1;
        end
        function r = pagetimes(o1, o2)
            %Realise le produit terme à terme de phaseurs de deux PhasorArray
            d=PhasorUnif(o1,o2);
            r = PhasorArray(pvalue(d{1}).*pvalue(d{2}));
        end
        function r = times(01,02)
        %TIMES Element-wise multiplication of two PhasorArray objects.
            R = TIMES(01,02) returns a PhasorArray containing the element-
wise
            multiplication of O1 and O2. Both inputs must be PhasorArray
objects
            (or unify to the same dimension).
            d=PhasorUnif(o1,o2);
            o1=d\{1\};
            o2=d\{2\};
            for ii = 1:size(o1,1)
            for jj = 1:size(o1,2)
                if ii==1 && jj==1
                rr=PhasorArrayTimes(o1(ii,jj,:),o2(ii,jj,:),reduce=false);
                r=zeros(size(o1,1), size(o1,2), size(rr,3));
                r(ii,jj,:)=rr;
                else
r(ii,jj,:)=PhasorArrayTimes(o1(ii,jj,:),o2(ii,jj,:),reduce=false);
                end
            end
            end
            r = PhasorArray(r);
        function r = mtimes(o1, o2)
            %MTIMES Overloads the matrix multiplication operator for
PhasorArray.
               R = MTIMES(01,02) performs the time-domain product of the two
                PhasorArray objects O1 and O2 (convolution of the 3D arrays
along the third dimension), returning a PhasorArray result.
            r = PhasorArray(PhasorArrayTimes(o1,o2));
        end
        function r = rdivide(o1, o2, varargin)
        % RDIVIDE Overloads the right array division operator for the
PhasorArray class.
        % RDIVIDE(A,B) returns a PhasorArray that represents A./B term by
term.
            for ii = 1:size(o2,1)
            for jj = 1: size(02,2)
                d2(ii,jj,:)=PhasorInv(o2(ii,jj,:),varargin{:});
            end
            end
            d2=PhasorArray(d2);
            d=PhasorUnif(o1,d2);
            r = PhasorArray(d{1}.*d{2});
```

```
function r = ldivide(o1,o2,varargin)
        % LDIVIDE Overloads the left array division operator for the
PhasorArray class.
            LDIVIDE(A,B) returns a PhasorArray that represents A.\B term by
term.
            for ii = 1:size(o1,1)
            for jj = 1:size(o1,2)
                d1(ii,jj,:)=PhasorInv(o1(ii,jj,:),varargin{:});
            end
            end
            d1=PhasorArray(d1);
            d=PhasorUnif(d1,o2);
            r = PhasorArray(d{1}.*d{2});
        end
        function r = pagerdivide(o1, o2)
            %Realise la r division terme à terme des Phaseurs de 2
            %PhasorArray
            d=PhasorUnif(o1,o2);
            r = PhasorArray(d{1}./d{2});
        end
        function r = pageldivide(01,02)
            %Realise la 1 division terme à terme des Phaseurs de 2
            %PhasorArray
            d=PhasorUnif(o1,o2);
            r = PhasorArray(d{1}.\d{2});
        function r = mrdivide(o1,o2,varargin)
            %MRDIVIDE Overloads the right matrix division operator for the
PhasorArray class.
            응
               MRDIVIDE(A,B) returns a PhasorArray that represents A(t)/
B(t), which is
                equivalent to A * inv(B) in the time domain.
               This function performs the time-domain matrix division of
the two
                PhasorArray objects 01 and 02.
            d2=PhasorInv(o2, varargin(:));
            r = PhasorArray(PhasorArrayTimes(o1,d2));
        function r = mldivide(o1, o2, varargin)
            %MLDIVIDE Overloads the left matrix division operator for the
PhasorArray class.
               MLDIVIDE(A,B) returns a PhasorArray that represents A(t)
\B(t), which is
                equivalent to inv(A) * B in the time domain.
                This function performs the time-domain matrix division of
the two
               PhasorArray objects 01 and 02.
            d1=PhasorInv(o1, varargin{:});
            r = PhasorArray(PhasorArrayTimes(d1,o2));
        end
        function r = pagepower(o1,m)
```

end

```
% pagepower Element-wise exponentiation of phasors.
           응
               r = pagepower(o1, m) computes `o1 .^ m` phasor-wise,
applying exponentiation
           응
               to each phasor coefficient individually.
           응
           용
              Inputs:
           응
              - ol: A PhasorArray.
               - m : A scalar or a PhasorArray of matching size.
              Behavior:
              - If `m` is a scalar, applies `.^m` to each individual
           응
phasor of `o1`.
               - If `m` is a PhasorArray, exponentiation is applied to
matching phasor coefficients.
               - Each frequency component of `o1` is exponentiated
separately.
              Notes:
              - This is **not** equivalent to exponentiating the time-
domain matrix \setminus ( A(t) \setminus).
               - Rarely useful in control or signal processing applications.
           9
               See also: power, mpower
           if isa(m,'PhasorArray')
               d=PhasorUnif(o1,m);
               m=d\{2\};
               r = PhasorArray((pvalue(d{1})).^(pvalue(m)));
           else
               r = PhasorArray(pvalue(o1).^m);
           end
        end
        function r = power(o1, m)
           % POWER Element-wise power of a periodic matrix A(t).
              r = power(o1, m) computes the element-wise power of each
matrix entry of a periodic matrix \setminus (A(t) \setminus),
               represented as a PhasorArray. This is equivalent to
computing \ (A(t)^{m}\ ) for each individual element of the matrix
              at each time instance in the time domain.
           응
           응
           응
              Inputs:
               - ol: A PhasorArray representing the periodic matrix \( A(t)
\).
           응
              - m: The exponent (scalar value) to apply element-wise.
           응
               Output:
               - r: A PhasorArray representing the matrix \( A(t)^{m} \)
computed element-wise.
              Behavior:
           % - The function applies the power operation **element-wise**
- This corresponds to raising each individual element \
```

```
( A \{ij\}(t) \) of the matrix to the power \( m \).
            응
                See also: PhasorPow, mpower, logm, expm
            r=o1;
            % element wise power
            for ii=1:size(o1,1)
                for jj=1:size(o1,2)
                    r{ii,jj}=o1{ii,jj}^m;
                end
            end
        end
        function r = mpower(o1, m)
            % MPOWER Matrix power of a periodic matrix A(t), computed for
integer exponents.
              r = mpower(01, m) computes the matrix power \( A(t)^m \) for
a periodic matrix \( A(t) \),
               represented as a PhasorArray. This operation is only valid
for **integer exponents** and is done in the time domain.
            % It computes the repeated matrix multiplication of \( A(t) \)
with itself \ (\ m\ ) times.
            응
            응
               Inputs:
               - ol: A PhasorArray representing the periodic matrix \((A(t))
\).
               - m: A positive or negative integer exponent representing
the power.
            응
            응
               Output:
               - r: A PhasorArray representing \( A(t)^m \), the matrix
raised to the \( m \)-th power.
              Behavior:
            % - For positive integer \( m \), it multiplies \( A(t) \)
with itself \( m \) times.
            % - For negative integer \( m \), it computes the inverse of \
(A(t)) raised to the power (|m|) and then multiplies.
               - For non integer \((m\), it calls PhasorPow, which perform
time domain matrix power and then IFFT.
            응
               Notes:
            % - This function applies matrix exponentiation, unlike
`power`, which operates element-wise.
            응
                See also: PhasorPow, power, logm, expm
            if mod(m, 1) == 0 \&\& m > 1
                prec=pvalue(o1);
                for ii=2:m
                    prec=PhasorArrayTimes(prec,o1);
            elseif mod(m, 1) == 0 \&\& m < 0
                d1=PhasorInv(o1);
                prec=d1;
                for ii=-2:-1:m
```

```
prec=PhasorArrayTimes(prec,d1);
                end
            else
                prec=PhasorPow(o1,m);
            end
            r=PhasorArray(prec);
        end
        function r = oplus(o1, o2)
            %OPLUS Kronecker sum of two PhasorArray objects.
                R = OPLUS(01,02) returns a PhasorArray representing the
Kronecker
            응
                sum of O1 and O2.
               The kronecker sum A(t) oplus B(t) is usually defined as
            양
                       A(t) otimes I + I otimes B(t)
                where I is the identity matrix of the same size as A(t) and
B(t)
            응
               and otimes is the usual kronecker product.
                see also KRON, TROPLUS, CTROPLUS
            r=PhasorArray(PhasorArrayOplus(o1,o2));
        end
        function r = troplus(01,02)
            %TROPLUS Kronecker sum of the transpose of two PhasorArray
objects.
                R = TROPLUS(01,02) returns a PhasorArray representing the
Kronecker
            응
               sum of the transpose of O1 and O2.
                see also KRON, OPLUS, CTROPLUS
            r=PhasorArray(PhasorArrayOplus(pagetranspose(pvalue(o1)),o2));
        function r = \text{ctroplus}(01,02)
            %CTROPLUS Kronecker sum of the conjugate transpose of two
PhasorArray objects.
               R = CTROPLUS(01,02) returns a PhasorArray representing the
Kronecker
               sum of the conjugate transpose of O1 and O2.
                see also KRON, OPLUS, TROPLUS
            r=PhasorArray(PhasorArrayOplus(pagectranspose(pvalue(o1)),o2));
        end
        function r = kron(o1,o2)
            %KRON Kronecker product of two PhasorArray objects.
                R = KRON(01,02) returns a PhasorArray representing the
Kronecker
            응
               product of O1 and O2.
                see also OPLUS, TROPLUS, CTROPLUS
            r=PhasorArray(PhasorArrayKron(o1,o2));
        function r = retro(01)
            %RETRO Reverse the time axis of the PhasorArray.
```

```
B = RETRO(A) returns a PhasorArray B such that B(t) = A(-t).
                This function flips the PhasorArray along the third
dimension,
                effectively reversing the time axis.
            r=PhasorArray(flip((pvalue(o1)),3));
        end
        function r = trretro(o1)
            %TRRETRO Reverse the time axis and transpose the PhasorArray.
                B = TRRETRO(A) returns a PhasorArray B such that B(t) = A(-1)
t).'
                This function flips the PhasorArray along the third
dimension,
                effectively reversing the time axis and transposing each
page.
            r=PhasorArray(flip(pagetranspose(pvalue(o1)),3));
        end
        function r = ctrretro(o1)
            %CTRRETRO Reverse the time axis and conjugate transpose the
PhasorArray.
                B = CTRRETRO(A) returns a PhasorArray B such that B(t) = A(-
t)'.
               This function flips the PhasorArray along the third
dimension.
                effectively reversing the time axis and conjugate
transposing each page.
            r=PhasorArray(flip(pagectranspose(pvalue(o1)),3));
        end
        function out = extract(o1,index,symmetric)
            %EXTRACT Extract specified phasors from a PhasorArray.
                OUT = EXTRACT(O1, INDEX) returns a PhasorArray containing
only the
                phasors specified by INDEX. All other phasors are set to
zero.
                OUT = EXTRACT(O1, INDEX, SYMMETRIC) allows for symmetric
extraction.
                If INDEX contains only positive integers and SYMMETRIC is
true, the
                corresponding negative phasors are also included. If INDEX
            응
contains
                negative integers, SYMMETRIC is forced to false.
            응
            응
                INPUTS:
                              - The PhasorArray object to extract from.
                    01
                              - A logical array matching the third dimension
                    INDEX
of the
                                PhasorArray or a vector of positive and
            응
negative integers.
                    SYMMETRIC - (Optional) A logical flag indicating whether
to include
            응
                                symmetric phasors. Defaults to true.
            응
                OUTPUT:
```

```
- A PhasorArray containing only the specified
phasors.
            응
                EXAMPLES:
            응
                    A = PhasorArray.random(3, 3, 5);
            응
                    B = A.extract([1, 3, 5]);
                    C = A.extract([1, 3], false);
            응
            응
                    D = A.extract([1, 3], true);
                See also trunc, reduce, neglect
            arguments
            01
            index
            symmetric=true
            end
            if islogical(index)
            if numel(index)~=size(o1,3)
                error('logical array size must match third dim of
phasorArray')
            end
            index=find(index);
            %find complement of index
            indexc=setdiff(1:size(o1,3),index);
            out = 01;
            out(:,:,indexc)=0;
            else
            if all(index>=0) && symmetric
                index=[index -index];
                maxIndex = max(max(index),o1.h);
                indexc=setdiff(-maxIndex:maxIndex,index);
                out = 01;
                out{:,:,indexc}=0;
            else
                out = 01;
                maxIndex = max(max(abs(index)),o1.h);
                indexc=setdiff(-maxIndex:maxIndex,index);
                out{:,:,indexc}=0;
            end
            end
        end
        function r = trunc(o1, m)
            %TRUNC Truncate the PhasorArray to a specified number of phasors.
                R = TRUNC(O1, M) returns a new PhasorArray that is truncated
to
                contain only M phasors. If M is not specified, the function
will
            응
                use a default value.
            응
            응
                INPUTS:
            응
                    ol - The original PhasorArray object to be truncated.
                    m - (Optional) The number of phasors to retain in the
truncated
```

OUT

```
PhasorArray. If not provided, a default value is
used.
               OUTPUT:
            응
                   r - The truncated PhasorArray object containing only M
phasors.
            용
            응
               EXAMPLE:
                   A = PhasorArray.random(3, 3, 5);
                   B = A.trunc(3);
            용
                   % B now contains only 3 phasors from the original
PhasorArray A.
              See also REDUCE, NEGLECT
           arguments
           01
           m = []
           r=PhasorArray(ReduceArray(o1,m));
        end
        function r = reduce(o1,htrunc,varg)
            %REDUCE Truncate or filter the PhasorArray based on harmonic
order or magnitude thresholds.
            % This function reduces a PhasorArray by either:
            % - Truncating harmonics beyond a given order (`htrunc`).
               - Filtering harmonics based on their magnitude
(`reduceMethod` and `reduceThreshold`).
              - Optionally controlling how the 0th harmonic is handled
(`excludeOPhasor`).
               - Applying a hard or soft thresholding strategy
(`hardThresholdPhasors`).
              Syntax:
              r = REDUCE(o1)
                  Reduces `o1` using the default relative thresholding
method,
                   preserving all harmonics unless their magnitudes are
below 1e-15.
            % r = REDUCE(o1, htrunc)
                   Truncates `o1` to harmonics of order ≤ `htrunc`.
            % r = REDUCE(o1, htrunc, 'reduceMethod', method,
'reduceThreshold', threshold,
                             'excludeOPhasor', exclude,
'hardThresholdPhasors', hardThreshold)
                   Applies a combination of truncation and magnitude
filtering based on additional parameters.
            응
            % Input Arguments:
            % - o1 (PhasorArray) : The PhasorArray object to be reduced.
               - htrunc (integer, optional) : The maximum harmonic order to
```

```
retain. Default is empty ([]),
                  meaning harmonics are only removed based on filtering
criteria.
           % Name-Value Pair Arguments:
           % - 'reduceMethod' (char, optional) : Defines the reduction
approach.
                   - 'absolute' : Removes phasors whose magnitude is below
the **absolute** threshold.
                   - 'relative' (default) : Removes phasors whose magnitude
is below a **relative** threshold.
              - 'reduceThreshold' (double, optional) : The threshold value
for filtering phasors.
           - If 'absolute', removes phasors with magnitude <</pre>
`reduceThreshold`.
                   - If 'relative', removes phasors with magnitude <
`reduceThreshold * max(magnitude)`.
                  - Default: 1e-15.
              - 'excludeOPhasor' (logical, optional) :
                  - true (default) : The Oth harmonic (DC component) is
always considered in max magnitude evaluation.
                  - false : The Oth harmonic is ignored when computing the
relative maximum magnitude.
                            This affects thresholding in 'relative' mode.
           9
           % - 'hardThresholdPhasors' (logical, optional) :
                  - false (default) : `o1` is truncated to order `m`,
meaning **only harmonics beyond order m**
                     are removed when all phasors of order > `m` fall below
the threshold.
                     Lower-order phasors (< `m`) are **never** removed,
even if below threshold.
                - true : Any phasor below the threshold is explicitly
**set to zero**. Then, trailing zero phasors
                  (i.e., those bearing no information) are removed to
optimize storage.
           % Output:
           % - r (PhasorArray) : The reduced PhasorArray after applying
truncation or filtering.
           용
              Example:
           % % Truncate PhasorArray to the 5th harmonic
              r = reduce(o1, 5);
              % Apply absolute thresholding without truncating harmonics
           % r = reduce(o1, [], 'reduceMethod', 'absolute',
'reduceThreshold', 1e-10, 'excludeOPhasor', false);
           % Apply soft thresholding with truncation logic
           % r = reduce(o1, [], 'hardThresholdPhasors', false,
'reduceMethod', 'relative', 'reduceThreshold', 1e-12);
```

```
응
              % Apply hard thresholding where all small phasors are
explicitly set to zero
              r = reduce(o1, [], 'hardThresholdPhasors', true,
            용
'reduceMethod', 'relative', 'reduceThreshold', 1e-12);
               See also: ReduceArray
            arguments
                01
                htrunc=[]
               varg.reduceMethod char {mustBeMember(varg.reduceMethod,
{'absolute', 'relative'})} = 'relative'
               varg.reduceThreshold {mustBeNumeric,mustBeReal} = 1e-15
               varg.excludeOPhasor (1,1) logical = true
                varq.hardThresholdPhasors=false
            end
r=PhasorArray(ReduceArray(o1, htrunc, reduceMethod=varg.reduceMethod, reduceThre
shold=varg.reduceThreshold,excludeOPhasor=varg.excludeOPhasor,hardThresholdPh
asors=varg.hardThresholdPhasors));
        end
        function r = neglect(o1, reduceThreshold, varg)
            %NEGLECT Set to zero all phasors below a given threshold in a
PhasorArray.
               This function filters out low-magnitude phasors in a
PhasorArray `o1`,
               setting them to zero based on a specified threshold and
reduction method.
               Syntax:
              r = NEGLECT(o1, reduceThreshold)
                   Sets to zero all phasors in `o1` with magnitude below
`reduceThreshold`,
            용
                   using the default 'relative' thresholding method.
              r = NEGLECT(o1, reduceThreshold, 'reduceMethod', method,
'excludeOPhasor', exclude, 'h', h)
            응
                   Applies phasor filtering with additional options.
              Input Arguments:
            % - o1 (PhasorArray) : The PhasorArray object whose small
phasors should be neglected.
              - reduceThreshold (double) : The magnitude threshold below
           응
which phasors are set to zero.
              Name-Value Pair Arguments:
              - 'reduceMethod' (char, optional) : Defines how the
threshold is applied.
                    - 'absolute' : A phasor is set to zero if its magnitude
is below `reduceThreshold`.
                   - 'relative' (default) : A phasor is set to zero if its
magnitude is below
                      `reduceThreshold * max(magnitude)` (where the
```

```
reference maximum can exclude the Oth phasor).
               - 'excludeOPhasor' (logical, optional) :
                  - false (default) : The 0th phasor (DC component) **is
included** when computing the reference magnitude.
                   - true : The Oth phasor is **ignored** when computing
          용
the reference magnitude for relative thresholding.
               - 'h' (integer, optional) :
           응
                  - If provided, reduces the PhasorArray to at most `h`
harmonics after applying the threshold.
                  - Similar to `reduce(h)`, but performed after zeroing
small phasors.
            응
           응
              Output:
           % - r (PhasorArray) : The filtered PhasorArray, with phasors
below the threshold set to zero.
              Example:
              % Set to zero all phasors with magnitude below 1e-15 using
default relative method
           % r = neglect(o1, 1e-15);
              % Apply absolute thresholding (removes all phasors with
magnitude < 1e-15)
              r = neglect(o1, 1e-15, 'reduceMethod', 'absolute');
           응
              % Use relative thresholding, ignoring the 0th phasor when
computing the max reference
           % r = neglect(o1, 1e-15, 'reduceMethod', 'relative',
'excludeOPhasor', true);
               % Apply thresholding and then truncate to at most 5 harmonics
              r = neglect(o1, 1e-12, 'h', 5);
               See also: REDUCE, TRUNC, ReduceArray
           arguments
               01
               reduceThreshold {mustBeNumeric,mustBeReal} = 1e-15
               varg.reduceMethod {mustBeMember(varg.reduceMethod,
{'absolute', 'relative'})} = 'relative'
               varg.excludeOPhasor (1,1) logical = false
               varq.h=[]
           end
           val=o1.value;
           h=o1.h;
           switch varg.reduceMethod
               case 'absolute'
                   val rel=val;
               otherwise
                    if varg.exclude0Phasor
                       ref=max(abs(val(:,:,[1:h, (h+2):(2*h+1)])),[],3);
%maximum harmonic on each coeef, excepting the phasor 0.
                    else
```

```
ref=max(abs(val),[],3); %maximum harmonic on each
coeef, excepting the phasor 0.
                    val rel=abs(val./ref);
            end
            val(abs(val rel)<reduceThreshold)=0;</pre>
            r = PhasorArray(val);
r.reduce(varg.h, "reduceMethod", varg.reduceMethod, "reduceThreshold", reduceThre
shold, "excludeOPhasor", varg.excludeOPhasor);
        end
        function r = flip(o1, dim)
            %FLIP the PhasorArray along the dim dimension
                - dim : dimension to flip
                - r = flip(ol,dim) flip the PhasorArray along the dim
dimension
            응
                See also FLIPLR, FLIPUD
            응
            9
                Inputs:
            응
                    o1 - PhasorArray object to be flipped
            응
                    dim - Dimension along which to flip the PhasorArray
            응
            응
            응

    r - Resulting PhasorArray after flipping

            응
            응
                Example:
            응
                    pa = PhasorArray(rand(3,3,3));
            응
                    pa flipped = flip(pa, 2);
            응
            응
               Note:
                    If dim is 3, a warning is issued indicating that the flip
                    along the third dimension produces M(-t).
            if dim==3
                warning("flip along third dimension, produces M(-t)")
            r=PhasorArray(flip(o1.value,dim));
        end
        function r = fliplr(01)
            %FLIPLR Flip the PhasorArray left to right
                r = FLIPLR(o1) flips the PhasorArray o1 in the left-right
direction.
            응
                The function returns a new PhasorArray r with the elements
flipped.
            응
                Example:
            응
                    pa = PhasorArray(rand(3,3,3));
            응
                    pa flipped = fliplr(pa);
            응
                See also FLIP, FLIPUD
            r=PhasorArray(fliplr(o1.value));
```

```
end
        function r = flipud(01)
            %FLIPUD Flip the PhasorArray up to down
                r = FLIPUD(o1) flips the PhasorArray o1 in the up-down
direction.
            응
                The function returns a new PhasorArray r with the elements
flipped.
               Example:
            응
                   pa = PhasorArray(rand(3,3,3));
            응
                   pa flipped = flipud(pa);
            응
            응
              See also FLIP, FLIPLR
            r=PhasorArray(flipud(o1.value));
        end
        function r = rot90 (o1, varargin)
            %ROT90 Rotate the PhasorArray by 90 degrees
                r = ROT90(o1) rotates the PhasorArray object o1 by 90 degrees
               counterclockwise and returns the result in r.
               r = ROT90(o1, k) rotates the PhasorArray object o1 by 90
degrees
            응
               counterclockwise k times.
            응
            응
               Input:
                    o1 - PhasorArray object to be rotated
            응
                    k - (Optional) Number of times to rotate by 90 degrees
            응
            응
               Output:
                    r - PhasorArray object after rotation
                see also FLIP, FLIPLR, FLIPUD
            r=PhasorArray(rot90(o1.value, varargin(:)));
        end
        function r = horzcat(o1, varargin)
            %HORZCAT Concatenate PhasorArray objects horizontally
                R = HORZCAT(O1, VARARGIN) concatenates the PhasorArray
object 01
                with additional PhasorArray objects provided in VARARGIN,
                horizontally, if their vertical dimensions are compatible.
            응
            응
               Input arguments:
               01 - The first PhasorArray object.
               VARARGIN - Additional PhasorArray objects to concatenate
with O1.
            응
               Output arguments:
               R - The resulting PhasorArray object after horizontal
concatenation.
```

```
응
                If only one input argument is provided, the function returns
01.
                If the third dimension sizes of O1 and the current object in
VARARGIN
                are not equal, the function uses PhasorUnif to unify them
before
            응
                concatenation.
            양
            응
                Example:
            응
                    r = horzcat(o1, o2, o3);
                See also: VERTCAT, PHASORUNIF, PVALUE
            if nargin==1
                r=o1;
                return
            end
            for ii=1:numel(varargin)
                o2=vararqin{ii};
                if size(o1,3) ~= size(o2,3)
                    d=PhasorUnif(o1,o2);
                    r = PhasorArray([pvalue(d{1}) pvalue(d{2})]);
                else
                    u1=pvalue(o1);
                    u2=pvalue(o2);
                    r = PhasorArray([u1, u2]);
                end
                o1=r;
            end
        end
        function r = vertcat(varargin)
            %VERTCAT Concatenate PhasorArray objects vertically
                R = VERTCAT(01, 02, ...) concatenates the PhasorArray object
01
                with additional PhasorArray objects provided in VARARGIN,
            응
                vertically, if their horizontal dimensions are compatible.
            응
                Input arguments:
                O1 - The first PhasorArray object.
                VARARGIN - Additional PhasorArray objects to concatenate
with 01.
            응
            응
                Output arguments:
                R - The resulting PhasorArray object after vertical
concatenation.
                If only one input argument is provided, the function returns
01.
                If the third dimension sizes of O1 and the current object in
VARARGIN
                are not equal, the function uses PhasorUnif to unify them
before
            응
                concatenation.
```

```
용
                    r = vertcat(o1, o2, o3);
            응
                See also: HORZCAT, PHASORUNIF, PVALUE
            응
            if nargin==1
            r=varargin{1};
            return
            end
            if nargin>=2
            r=cat(1, varargin(:));
            return
            else
            o1=vararqin{1};
            o2=varargin{2};
            end
            if size(o1,3)~=size(o2,3)
            d=PhasorUnif(o1,o2);
            r = PhasorArray([pvalue(d{1}); pvalue(d{2})]);
            else
            u1=pvalue(o1);
            u2=pvalue(o2);
            r = PhasorArray([u1; u2]);
            end
        end
        function r = blkdiag(o1)
            %BLKDIAG of phasorarray
            % - r = blkdiag(o1,o2,...) : block diagonal of phasorArray
            응
               See also HORZCAT, VERTCAT
            응
               This function creates a block diagonal matrix from the input
               phasor arrays. It takes multiple phasor arrays as input and
               constructs a block diagonal matrix where each block
corresponds
            응
               to one of the input phasor arrays.
            응
            응
                Input:
            응
                    ol - Repeating argument list of phasor arrays
            응
            응
               Output:
            용
                    r - Resulting PhasorArray object containing the block
            응
                        diagonal matrix
            응
            응
               Example:
            응
                   pa1 = PhasorArray(rand(3,3,2));
            응
                   pa2 = PhasorArray(rand(2,2,2));
            응
                    result = blkdiag(pa1, pa2);
            응
                This example creates a block diagonal matrix from two phasor
                arrays pal and pa2.
            arguments (Repeating)
```

응

Example:

```
01
            end
            b1=PhasorUnif(o1{:});
            h=size(b1{1},3);
            out=cell(h,1);
            for hi=1:h
                argi=cell(numel(b1),1);
                for b1i=1:numel(b1)
                    bb1i=pvalue(b1{b1i});
                    argi{bli}=bbli(:,:,hi);
                end
                out{hi}=blkdiag(argi{:});
            end
            out=cat(3,out{:});
            r=PhasorArray(out);
        end
        function r = repmat(o1, M, N)
            %REPMAT Replicate and tile a PhasorArray.
               r = REPMAT(o1,M,N) replicates and tiles the PhasorArray o1 M
times along the first dimension and N times along the second dimension.
               r = REPMAT(o1, M) replicates and tiles the PhasorArray o1 M
times along the first and second dimensions.
            % r = REPMAT(o1,M,N,L) replicates and tiles the PhasorArray o1
M times along the first dimension, N times along the second dimension, and L
times along the third dimension.
            응
            응
                Inputs:
                    ol - The PhasorArray object to be replicated.
                    \mbox{\tt M} - Number of times to replicate along the first
dimension.
                    N - (Optional) Number of times to replicate along the
second dimension.
            응
                Outputs:
            응
                    r - The resulting PhasorArray after replication.
            응
            응
                Example:
            응
                    pa = PhasorArray(rand(3,3,3));
            응
                    pa replicated = repmat(pa, 2, 3);
            응
                See also REPMAT, RESHAPE
            arguments
            01
            Μ
            arguments (Repeating)
            end
            switch nargin
            case 1
                r = 01;
            case 2
                switch numel(M)
```

```
case 1
                    N=M;
                    L=1;
                    M = [M N L];
                case 2
                    N=M(2);
                    M=M(1);
                    L=1;
                    M = [M N L];
                case 3
                    warning("You shouldn't specify a third dimension for
repmat with phasorarray, proceed with caution")
                otherwise
                    warning("You shouldn't specify more than 2 dimensions
for repmat with phasorarray, proceed with caution")
                end
                r = PhasorArray(repmat(o1.value,M));
                return
            case 3
                assert(numel(M) ==1)
                N=N\{1\};
                assert(numel(N) ==1)
                L=1;
                r = PhasorArray(repmat(o1.value, M, N, L));
                return
            otherwise
                warning("You shouldn't specify more than 2 dimensions for
repmat with phasorarray, proceed with caution")
                r = PhasorArray(repmat(o1.value,M,N(:)));
                return
            end
        end
        function r = reshape(o1, varargin)
            % reshape Reshape a PhasorArray while preserving the phasor
dimension.
            응
                r = reshape(o1, M, N)
                                         - Reshapes `o1` into an MxN
PhasorArray,
                                             preserving the third (phasor)
dimension.
                r = reshape(o1, M, N, L) - Reshapes `o1` into an MxNxL
PhasorArray.
            응
                Notes:
                - If only `M, N` are provided, `L` is automatically set to
`size(o1,3)`.
                - The third dimension (phasor order) should remain unchanged
in most cases.
                See also: repmat, reshape
            if numel(varargin) == 2
                varargin{3}=o1.size(3);
            end
```

```
r=PhasorArray(reshape(pvalue(o1), varargin{:}));
        end
        function r = permute(o1, varargin)
            % permute Reorder dimensions of a PhasorArray.
                r = permute(o1, order) - Rearranges the dimensions of `o1`
according to `order`.
            용
               r = permute(o1)
                                        - Swaps the first and second
dimensions (acts as transpose).
            응
               Notes:
               - By default, swaps dimensions `[1 2]`, equivalent to a
transpose.
               - The third dimension (phasor order) is **preserved** unless
explicitly modified.
               - Usually applied to **matrix-like** PhasorArrays, not 3D
arrays.
              See also: transpose, permute
            if nargin==1
                varargin{1}=[2 1];
            end
            if numel(varargin{1}) == 2
                varargin{1}=[varargin{1} 3];
            end
            r=PhasorArray(permute(pvalue(o1), varargin(:)));
        end
        function r = sub(o1, n1, n2)
            % sub Extract specific elements from a PhasorArray based on
indices.
                r = sub(o1, n1, n2) extracts the phasor array component
            응
                `Phasor3D(n1, n2, :)`, equivalent to `A{n1, n2}`.
            응
            응
                Inputs:
                   ol - The PhasorArray object to extract from.
                   n1 - Row indices or logical mask (can be `:` for all
rows).
                    n2 - (Optional) Column indices (default: 1 if o1 is a
column vector).
            응
                Output:
                    r - Extracted PhasorArray elements with `reduce=false`
            응
to
            응
                          preserve the structure.
            응
                Behavior:
                - If `n1` is a logical array, it is converted to indices.
               - If `o1` is a column vector (`size(o1,2)==1`), `n2`
defaults to 1.
               - If `n1` and `n2` are omitted, the function reshapes `r` to
match
                  size(n1,1) \times size(n1,2).
            응
```

```
See also: PhasorArray, reduce, logical indexing
            if nargin<3 %cas ou on a en apparence que une seule entrée de
coordonnée
                % cas 1 c'est une matrice d'entier
                % cas 2 c'est une matrice de logical, et il faut verif
                % qu'elle fait la bonne taille
                if ~strcmp(n1,':')%on a une matrice en input
                     if islogical(n1)
                         if numelt(o1) == numel(n1)
                             n1=find(logical(n1));
                         else
                             error('logical array size %d x %d must match
first two dim of phasorArray %d x
%d', size (n1,1), size (n1,2), size (o1,1), size (o1,1))
                         end
                     end
                end
                if size(o1,2)==1 % c'est déjà un phasorArray colonne, normal
de n'avoir eu qu'un indice
                    n2=1;
                else
                    o1=vect(o1); % on le rend colonne, ce sera plus simple à
manipuler
                    n2=1;
                end
            end
            r=PhasorArray(o1(n1,n2,:),reduce=false);
            if nargin == 1
            r = reshape(r, size(n1, 1), size(n1, 2));
            end
        end
        function r = vect(o1)
           %VECT transform phasorArray matrix to column vector (col operator)
                This function takes a PhasorArray object and transforms it
into
                a column vector by stacking the columns of the input matrix.
            응
            응
                Syntax:
            응
                    r = vect(o1)
            응
            용
                Input:
            응
                    o1 - PhasorArray object to be transformed.
            응
            응
                Output:
                    r - Column vector (PhasorArray) obtained by stacking the
            응
            응
                         columns of the input matrix.
            응
            응
                Example:
            응
                    % Assuming o1 is a PhasorArray object
            응
                    r = vect(o1);
```

```
See also: reshape, pvalue
            r1=reshape(pvalue(o1),[],1,size(o1,3));
            r=PhasorArray(r1, reduce=false);
        end
        function r = pad(o1,delta h)
            %PAD the PhasorArray with delta h phasor
               - r = pad(o1,delta h) : pad the PhasorArray o1 with delta h
0 phasor
                -r = pad(o1, [delta h1 delta h2 delta h3]) : pad the
PhasorArray o1 with delta h1 0 phasor along the first dimension, delta h2 0
phasor along the second dimension and delta h3 0 phasor along the third
dimension
            %pad (add zeros) to phasor array
            %pad(A, h) add h phasor to the phasor array
            %pad(A, [h1 h2 h3]) pad A in each direction
            r=PhasorArrayPad(o1,delta h);
        end
        function r = ctranspose(o1)
            %CTRANSPOSE overloading for PhasorArray
            CTRANSPOSE(A) : A' = A(t)'
            %transpose avec '
            r = mctranspose(o1);
        end
        function r = transpose(o1)
            %TRANSPOSE overloading for PhasorArray
            TRANSPOSE(A) : A.' = A(t).'
            %transpose avec .'
            r = mtranspose(o1);
        function r = pagectranspose(o1)
            %PAGECTRANSPOSE overloading for PhasorArray
            %PAGECTRANSPOSE(A) apply ctranspose to each page of A resulting
in A(-t)'
            r = PhasorArray(pagectranspose(pvalue(o1)));
        end
        function r = pagetranspose(o1)
            %PAGETRANSPOSE overloading for PhasorArray
            %PAGETRANSPOSE(A) apply transpose to each page of A resulting in
A(t).'
            % transpose appliqué à chacun des phaseurs
            r = PhasorArray(pagetranspose(pvalue(o1)));
        end
        function r = mtranspose(o1)
            % transpose appliqué à la matrice temporelle
            r=pagetranspose(o1);
        end
        function r = mctranspose(o1)
            % ctranspose appliqué à la matrice temporelle
            r=PhasorArray(flip(pagectranspose(o1.value),3));
        end
        function r = pmax(o1, o2, varg)
```

```
% pmax Compute elementwise or phasorwise max between two
PhasorArray objects.
                r = pmax(o1, o2, 'method', method) computes the maximum
            응
between two
                PhasorArray objects `o1` and `o2` according to the specified
method.
            응
                Inputs:
                   ol, o2 - PhasorArray objects to compare.
                    method - (Optional) Specifies the max computation
method:
                                * 'elementwise' (default): Max computed based
on norm.
                               * 'phasorwise': Max computed independently
for each phasor.
               Output:
                            - The resulting PhasorArray after applying the
                   r
max operation.
            응
               Notes:
               - In 'elementwise' mode, each phasor's norm is computed
first, and the
                 element-wise max is applied to select the dominant phasor.
            응
               - In 'phasorwise' mode, each corresponding phasor component
is compared directly.
                - Both inputs are unified in size using `PhasorUnif` before
processing.
                See also: PhasorUnif, pvalue, norm
            arguments
                01
                02
                varg.method char {mustBeMember(varg.method,
{'elementwise', 'phasorwise'})} = 'elementwise'
            end
            if strcmp(varg.method, 'elementwise')
                d=PhasorUnif(o1,o2);
                o1 = d\{1\};
                o2 = d\{2\};
                % norm of each phasor (along the third dim)
                n1=sum(o1.value .* conj(o1.value),3);
                n2=sum(o2.value .* conj(o2.value),3);
                %max of the norm
                n=\max(n1,n2);
                %elementwise max
                r=PhasorArray(pvalue(o1).*(n1==n) + pvalue(o2).*(n1\sim=n));
            else
                d=PhasorUnif(o1,o2);
                r = PhasorArray(max(pvalue(d{1}), pvalue(d{2})));
```

```
end
        end
        function varargout=phas(o1,h)
            % phas Extract the phasor of order `h` from a PhasorArray.
                phas(o1, h) returns the phasor of order `h` from the
PhasorArray `o1`.
            양
                Inputs:
                    ol - The PhasorArray object.
                    h - The order of the phasor to extract.
                         - If `h` is a scalar, returns a single phasor slice.
            응
                         - If `h` is a vector, returns multiple slices at
the specified orders.
            용
               Outputs:
            응
                   - A 3D array corresponding to the phasor(s) of order `h`.
            90
               Behavior:
               - If `abs(h) <= o1.h`, extracts the corresponding phasor.
               - If `h` is **out of range**, pads `o1` to match the highest
requested order.
               - If `h` is **beyond padding limits**, returns a zero matrix.
            응
               See also: pad, reduce, PhasorArray
            if isscalar(h)
                if abs(h)<=01.h
                    varargout{1}=o1(:,:,(end+1)/2+h);
                else
                    varargout{1}=zeros(size(o1,[1 2]));
                end
            else
                if \max(abs(h)) >= 01.h
                    o1=o1.pad( max(abs(h))-o1.h);
                varargout{1}=o1(:,:,(end+1)/2+h);
            end
        end
        function obj =phasAssign(obj,h,varargin)
            % phasAssign Assign values to specific phasor components in a
PhasorArray.
                obj = phasAssign(obj, h, values) returns a copy of the
PhasorArray
                with the phasor at index h replaced by the provided values.
            응
            응
                Inputs:
            응
                           - The PhasorArray object to modify.
                    obj
                           - The index (or indices) of the phasor(s) to be
assigned.
                   values - The new values to be assigned to the specified
```

```
phasors.
            응
            응
               Outputs:
                    obj
                          - A modified PhasorArray with updated phasor
values.
               Notes:
            % - If h exceeds the current order of the PhasorArray, it is
automatically expanded.
               - This method does not modify the original object but
returns a modified copy.
           % - The provided values must be compatible in size with the
existing array dimensions.
            응
            응
                See also: phas, sub, reduce, trunc
            obj{:,:,h}=varargin{:};
        end
        function r = lt(o1,o2)
            r= (pvalue(o1)<pvalue(o2));
        end
        function r = gt(o1,o2)
            r= (pvalue(o1)>pvalue(o2));
        end
        function r = le(o1, o2)
            r= (pvalue(o1) <=pvalue(o2));
        end
        function r = ge(o1, o2)
            r= (pvalue(o1)>=pvalue(o2));
        function r = ne(o1,o2)
            r= (pvalue(o1)~=pvalue(o2));
        end
        function r = eq(01,02)
            r= (pvalue(o1) == pvalue(o2));
        end
        function r = and(o1,o2)
            r= (pvalue(o1) & pvalue(o2));
        end
        function r = or(o1, o2)
            r= (pvalue(o1) | pvalue(o2));
        end
        function r = not(o1)
            r= not(pvalue(o1));
        end
        function r = double(o1)
            %DOUBLE Convert PhasorArray to double precision.
               r = DOUBLE(o1) converts the PhasorArray object o1 to a
double precision array.
            r= double(pvalue(o1));
        end
```

```
function o2 = repeat(o1,m)
            %REPEAT Repeat the PhasorArray object over its period by a
factor m
            응
            응
                o2 = repeat(o1, m) repeats the PhasorArray object o1 over its
                period by a factor m. If m is negative, the signal is
reversed
            응
                in time, effectively dividing the period by m.
            응
            응
                Inputs:
            응
                    ol - The PhasorArray object to be repeated.
                    m - An integer factor by which to repeat the period.
            응
            응
                         Default value is 2. If m is negative, the signal is
                         reversed in time.
            응
            응
            응
               Outputs:
                    o2 - The resulting PhasorArray object after repeating the
            응
            응
                         period by a factor m.
            응
            9
               Example:
                    o1 = PhasorArray(...); % Create a PhasorArray object
                    o2 = repeat(o1, 3); % Repeat the period of o1 by a
factor of 3
               See also: PhasorArray.zeros, retro
            arguments
                01
                m \{ mustBeInteger(m) \} = 2
            end
            h=o1.h;
            h new=h*abs(m);
            o2 = PhasorArray.zeros(size(o1,1), size(o1,2),2*h new+1);
            o2(:,:,1:abs(m):end)=o1.value;
            if m<0
                o2=retro(o2);
            end
        end
        function r = antiD(o1,T)
            %ANTI-D Compute the anti-derivative (primitive) of non-zero
phasors of a PhasorArray.
               r = ANTI-D(o1, T) computes the anti-derivative of the
            용
PhasorArray o1
               with respect to the period T. The function ignores the 0-th
phasor to
               produce a periodic output.
            응
            응
            응
               INPUTS:
                    o1 - The PhasorArray object to be anti-differentiated.
                    T - (Optional) The period of the PhasorArray. Defaults
to 2*pi.
```

```
응
               OUTPUT:
                    r - The resulting PhasorArray after computing the anti-
derivative.
            응
                EXAMPLE:
            응
                 A = PhasorArray.random(3, 3, 5);
            응
                    B = A.antiD(2*pi);
            응
            응
                NOTE:
                    The function divides each non-zero phasor by (1i * k * 2
* pi / T),
                   where k is the phasor index. The 0-th phasor is set to
zero.
            arguments
            01
            T=2*pi
            end
            o1=o1.value;
            [\sim, \sim, n3] = size(o1);
            for ii=1:n3
            k=-(n3-1)/2+ii-1;
            if k \sim = 0
            o1(:,:,ii)=o1(:,:,ii)/(1i*k*2*pi/T);
            else
                o1(:,:,ii)=0;
            end
            end
            r=PhasorArray(o1);
        end
        function r = d(o1,T)
            % d - Derive phasor array with respect to a given period
            % Syntax: r = d(o1, T)
            용
            % Inputs:
               o1 - The input phasor array object
                 T - The period with respect to which the phasor array is
derived (default is 2*pi)
            응
            % Outputs:
                r - The resulting phasor array after derivation
            % Example:
                 phasorArray = PhasorArray(someValues);
                 result = phasorArray.d(phasorArray, 2*pi);
            arguments
                01
                T=2*pi
            end
            %[\sim, \sim, n3] = size(o1);
            try
                K=permute((-o1.h):(o1.h),[1 3 2])*1i*2*pi/T;
                oo=bsxfun(@times,o1.value,K);
```

```
r=PhasorArray(oo);
            catch e
                % We need to manually perform the operation intended for
each "slice" of P.
                for idx = 1:size(o1, 3)
                    o1(:,:,idx) = o1(:,:,idx) * K(idx);
                r = 01;
            end
        end
        function r = PhaseShift(o1,angle)
            % PhaseShift Apply a phase shift to a PhasorArray.
            % r = PhaseShift(o1, angle) shifts the phase of the
PhasorArray `o1`
               by `angle` radians.
            용
            응
            용
               Inputs:
                         - The PhasorArray object.
                    01
                    angle - The phase shift value(s). It can be:
                           - A scalar (applies to all elements of `o1`).
                            - A row vector (broadcasts across `o1` columns
if `o1` is a column).
                           - A column vector (broadcasts across `o1` rows
if `o1` is a row).
                           - A matrix (applies element-wise if `o1` is
scalar).
            응
            % Output:
                   r - The phase-shifted PhasorArray.
            % Behavior:
              - If `o1` is a scalar, `angle` can be a **matrix**, applying
element-wise shifts.
              - If `o1` is a **row vector**, `angle` must be a column
vector (broadcasted).
           9
              - If `o1` is a **column vector**, `angle` must be a row
vector (broadcasted).
              - If `ol` is a **matrix**, `angle` must be scalar.
            응
            응
              Errors:
               - If `o1` and `angle` are both row or column vectors, the
function raises an error.
              - If `o1` is a matrix, `angle` must be a scalar.
                See also: dephase, PhasorArray
           if numel(angle)>1
                if isrow(o1) && iscolumn(angle)
                    n1=numel(angle);
                    %n2=o1.size(2);
                    r=repmat(o1, n1, 1);
                    for angli=1:numel(angle)
                        angle(angli);
```

```
r{angli,:}=o1.PhaseShift(angle(angli));
                     end
                    return
                elseif iscolumn(o1) && isrow(angle)
                     %n1=o1.size(1);
                    n2=numel(angle);
                    r=repmat(o1,1,n2);
                     for angli=1:numel(angle)
                         r{:,angli}=o1.PhaseShift(angle(angli));
                     return
                elseif isscalar(o1) && ismatrix(angle)
                     [n1, n2] = size (angle, [1, 2]);
                     r=repmat(o1,n1,n2);
                     for angli=1:size(angle,1)
                         for anglj=1:size(angle,2)
                             r{angli,anglj}=o1.PhaseShift(angle(angli,anglj));
                     end
                     return
                elseif iscolumn(o1) && iscolumn(angle)
                     error('if o1 is a column, angle must be row and vice
versa')
                elseif isrow(o1) && isrow(angle)
                    error('if o1 is a row, angle must be colmun and vice
versa')
                else
                    error('angle must be a scalar input for matricial
PhasorArray')
                end
            else
                r=dephase(o1, angle);
            end
        end
        function [oInv, oInvt, norm err, norm ref] = inv(o1, varargin)
            %INV Compute the phasor representation of the pointwise inverse
of A(t).
                This function computes the **pointwise inverse** of the time-
domain realization
                of the periodic matrix A(t) and then reconstructs its phasor
representation.
            응
                The inversion process follows these steps:
                1. Compute the **time-domain representation** A(t) using an
**IFFT**.
            응
                2. Perform **pointwise inversion** A<sup>-1</sup>(t).
            응
                3. Reconstruct phasors by applying an **FFT** to A^{-1}(t).
            응
            응
                Syntax:
                r = INV(o1)
            응
                    Computes the phasors of the pointwise inverse of `o1`
```

```
using default parameters.
                [r, At] = INV(01)
                    Returns the time-domain realization A^{-1}(t) alongside the
phasors.
            응
               [r, At, norm err, norm ref] = INV(01)
                    Also returns reconstruction error metrics.
            응
               Input Arguments:
               - o1 (PhasorArray) : The PhasorArray object representing
A(t).
            응
               Output Arguments:
               - oInv (PhasorArray): The phasor representation of A^{-1}(t).
               - oInvt (array, optional) : The time-domain realization of A^{-1}
(t).
            용
               - norm err (double, optional) : Reconstruction error ||
Ainv ph - Ainv t \mid \mid F.
               - norm ref (double, optional) : Reference norm ||Ainv t|| F.
            응
            응
            응
               Example:
               % Compute the phasors of the inverse of a given PhasorArray A
               r = inv(A);
               % Compute both the phasors and time-domain representation
            응
            응
               [r, At] = inv(A);
            응
            응
               % Compute full outputs including reconstruction errors
                [r, At, err, ref] = inv(A);
               See also: DET, REDUCE.
            % Call PhasorInv with only the required number of outputs
            if nargout == 1
                Ainvph = PhasorInv(o1, varargin{:});
                oInv = PhasorArray(Ainvph);
            elseif nargout == 2
                [Ainvph, oInvt] = PhasorInv(o1, varargin{:});
                oInv = PhasorArray(Ainvph);
                [Ainvph, oInvt, norm err, norm ref] = PhasorInv(o1,
varargin(:));
                oInv = PhasorArray(Ainvph);
            end
        end
        function [PhDet, det t] = det(o1, varg)
            %DET Compute the pointwise determinant of A(t) and reconstruct
its phasors.
               This function computes the determinant of the **time-domain
realization ** A(t),
            % then reconstructs its phasor representation via Fourier
Transform.
```

```
응
              The determinant is computed **pointwise at each time step**,
meaning:
              - If A(t) is periodic, det(A(t)) is computed over time.
               - The final result represents the phasor decomposition of
det(A(t)).
            응
            응
               Syntax:
               [PhDet, det t] = DET(o1)
                   Computes the phasors of det(A(t)) using default settings.
           용
           % [PhDet, det t] = DET(o1, 'nT', nT, 'T', T, 'm', m,
                                         'plot', plotFlag, 'autoTrunc',
autoTrunc)
                   Computes det(A(t)) with additional control over
truncation and plotting.
            % Input Arguments:
              - ol (PhasorArray) : The PhasorArray object whose
determinant is computed.
           용
              Name-Value Pair Arguments:
            % - 'nT' (integer, optional) : Number of periods used in the
time-domain evaluation. Default: 1.
           % - 'T' (double, optional) : The period used for simulation.
Default: 1.
            % - 'm' (integer, optional) :
                  - Power of two controlling time-domain discretization.
                  - Can be set to [] for automatic selection based on the
number of phasors.
           % - 'plot' (logical, optional) : If true, plots det(A(t))
after computation. Default: false.
            % - 'autoTrunc' (logical, optional) :
                  - true : Uses the derivative of phasors to
**automatically detect**
                     the significant number of phasors.
                   - false (default) : Uses a fixed threshold-based
reduction method.
            % - If 'autoTrunc' is false, the following options apply:
                   - 'reduceThreshold' (double, optional) : The threshold
for reducing phasors. Default: 1e-20.
                  - 'reduceMethod' (char, optional) : Reduction strategy.
                       - 'absolute' : Remove phasors with magnitude <
reduceThreshold.
                       - 'relative' (default) : Remove phasors with
magnitude < max(magnitude) * reduceThreshold.</pre>
              Output Arguments:
               - PhDet (PhasorArray) : The phasor representation of
det(A(t)).
              - det t (array) : The time-domain realization of det(A(t)).
              Algorithm:
```

```
1. Compute A(t) in the time domain using an **IFFT**.
                2. Perform **pointwise determinant computation** det(A(t)).
                3. Reconstruct phasors by applying an **FFT** to det(A(t)).
            응
            응
                Example:
                % Compute the phasors of det(A(t)) using default settings
            용
               [PhDet, det t] = det(A);
            응
               % Compute det(A(t)) over 2 periods with auto truncation
               [PhDet, det t] = det(A, 'nT', 2, 'autoTrunc', true);
            9
               % Compute det(A(t)) with manual truncation and thresholding
               [PhDet, det t] = det(A, 'reduceThreshold', 1e-15,
'reduceMethod', 'absolute');
            응
              See also: INV, REDUCE, FFT, IFFT.
            arguments
            01
            varg.nT=1
            varg.T=1
            varq.m=[]
            varg.plot=false
            varg.reduceThreshold = 1e-20
            varg.reduceMethod = 'relative'
            varg.autoTrunc = false
            end
            C=namedargs2cell(varg);
            [PhDet, det t] = PhasorDet(o1, C(:));
            PhDet=PhasorArray(PhDet);
        end
        function r = diag(o1, K)
            %DIAG Extract or construct a diagonal PhasorArray.
               This function operates similarly to MATLAB's `diag`:
               - If `o1` is a **vector**, it constructs a **diagonal
PhasorArray**.
               - If `ol` is a **square matrix**, it extracts its diagonal
as a **vector**.
            응
               Syntax:
            응
               r = DIAG(A)
                    Constructs a diagonal PhasorArray from a vector `A`.
               r = DIAG(A, K)
            응
                   Constructs a diagonal PhasorArray, with `K` as the
diagonal offset.
            응
              r = DIAG(A)
                    Extracts the diagonal as a PhasorArray vector when `A`
is square.
```

```
응
                Input Arguments:
            응
                - o1 (PhasorArray) : Input PhasorArray, either a vector or a
square matrix.
                - K (integer, optional) : Offset for the diagonal placement.
Default: 0.
            응
            응
                Output:
                - r (PhasorArray) :
                    - If `o1` is a vector, `r` is a **diagonal PhasorArray**.
            응
                    - If `ol` is square, `r` is the **extracted diagonal**.
            응
            응
                Example:
            응
                % Create a diagonal PhasorArray from a vector
            응
                A = PhasorArray(rand(5,1));
            응
                R = diag(A);
            응
            응
                % Extract the diagonal from a square PhasorArray
                B = PhasorArray(rand(4,4));
                d = diag(B);
            응
            응
                % Construct a diagonal PhasorArray with an offset
            응
                C = diag(A, 1);
               See also: TRACE.
            arguments
                01
                K=0
            end
            if isscalar(o1) && K==0
                r = 01;
                return
            end
            if isvector(o1)
                if ~iscolumn(o1)
                    01=01.';
                end
                if isa(o1.value, "sdpvar")
                    r = PhasorArray(diag(o1));
                    return
                end
                if isa(o1.value, "ndsdpvar") || isa(o1.value, "sym")
                    for ii = 1:size(o1,3)
                        r{ii} = PhasorArray(diag(o1(:,:,ii)));
                    r = cat(3, r\{:\});
                    return
                end
                r = PhasorArray.zeros(numelt(o1)+abs(K));
```

```
I = logical(diag(ones(numelt(o1),1),K));
                r{I}=r{I}+o1;
                return
            end
            assert(issquare(o1), "diag function: PhasorArray must represent
a square matrix");
            n=size(o1,1);
            I = (1:n) * (n+1) -n;
            r=o1{I};
        end
        function r = trace(o1, type)
            \mbox{\ensuremath{\$TRACE}} Compute the phasor representation of the trace of A(t) .
               This function computes the **trace** of a PhasorArray A,
which is the sum
              of its diagonal elements. It supports two modes:
               - 'phasor' (default) : Returns the **full** phasor
representation of the trace.
               - '0'
                                    : Returns only the **DC (0th order) **
            응
component of the trace.
            용
            응
               Syntax:
               r = TRACE(A)
                    Computes the phasor representation of the trace of `A`.
            응
            응
               r = TRACE(A, 'phasor')
            응
                    Same as above; returns the full phasor trace.
            응
            00
               r = TRACE(A, '0')
                    Returns only the DC component of the trace.
            응
               Input Arguments:
            % - o1 (PhasorArray) : A **square PhasorArray**.
            % - type (char, optional) :
                   - 'phasor' (default) : Returns the **full phasor
            응
representation ** of trace(A).
                  - '0': Returns only the **DC component** (0th phasor)
of trace(A).
              Output:
            % - r (PhasorArray or numeric) :
            응
                    - If 'phasor', `r` is a **PhasorArray** containing the
full phasor trace.
                   - If '0', `r` is a **scalar numeric value** (the DC
            용
component of trace(A)).
            응
               Example:
               % Compute the full phasor trace of a PhasorArray
            % A = PhasorArray(rand(3,3,11));
               tr phasor = trace(A);
            응
               % Compute only the DC component of the trace
               tr DC = trace(A, '0');
```

```
See also: DIAG.
            arguments
                01
                type char {mustBeMember(type, {'phasor', '0'})} = 'phasor'
            end
            r=sum(diag(o1));
            if strcmp(type,'0')
                %return the DC value of the trace
                r=r{:,:,0};
            end
        end
        function r = numelt(o1)
            %NUMELT Compute the number of elements in the first two
dimensions of A.
                This function returns `size(A,1) * size(A,2)`, which
represents
               the number of matrix elements per time instant.
            응
            응
               Syntax:
            응
               r = NUMELT(A)
            응
                    Computes the total number of elements in A(t).
            응
            응
                Input:
            응
               - o1 (PhasorArray) : A PhasorArray object.
            응
            응
                Output:
            응
                - r (integer) : Number of elements in `A(t)`.
               Example:
                % Compute the number of elements in a 4x5 PhasorArray
                A = PhasorArray(rand(4,5,10));
                e = numelt(A); % Returns 4*5 = 20.
            r=size(o1,1)*size(o1,2);
        end
        function r = h(o1)
            %H Compute the maximal phasor order stored in A.
            응
                This function extracts the **highest phasor order** present
in a PhasorArray.
            응
            응
                Syntax:
                r = H(A)
            응
                    Computes the highest stored harmonic order.
            응
            응
                Input:
            응
                - o1 (PhasorArray) : A PhasorArray object.
            응
            응
                Output:
            응
                - r (integer) : The highest stored harmonic order in A.
            응
                Example:
                % Get the maximum harmonic order in a PhasorArray
```

```
A = PhasorArray(rand(4,4,11));
                max h = h(A); % Returns (11-1)/2 = 5.
                See also: DIM.
            r = (size(o1, 3) - 1)/2;
        end
        function r = dim(o1)
            %DIM Compute the number of elements in A(t).
                This function returns `size(A,1) * size(A,2)`, which
represents
            응
               the number of matrix elements per time instant.
            응
            응
               Syntax:
            응
                r = DIM(A)
            응
                    Computes the total number of elements in A(t).
            응
            양
                Input:
            응
               - o1 (PhasorArray) : A PhasorArray object.
            응
               Output:
            응
                - r (integer) : Number of elements in `A(t)`.
            응
                Example:
            응
               % Compute the number of elements in a 4x5 PhasorArray
                A = PhasorArray(rand(4,5,10));
                e = dim(A); % Returns 4*5 = 20.
                See also: NUMELT.
            r=size(o1,1)*size(o1,2);
        end
        function r = BT(o1, m)
            %BT Construct a Block Toeplitz matrix from a PhasorArray.
                This function builds a **Block Toeplitz matrix** of order
`m` from the given
                PhasorArray `A`. A **Block Toeplitz matrix** is a **block-
structured matrix**
                where **entire blocks** (submatrices) are repeated along the
diagonals.
            용
                Specifically, the matrix `BT(A, m)` is given by:
                    B = sum \{k=-2m\}^{2m} \ (I^k \otimes A k),
                where:
                - `I^k` is the **diagonal selector matrix**, which has ones
on its `k`-th diagonal.
               - `⊗` denotes the **Kronecker product**.
                - `A k` are the Fourier components (harmonics) of `A`.
            응
                This matrix represents the **finite-dimensional truncation**
of an
               **infinite-dimensional Block Toeplitz operator** acting on
```

```
\ell^2, where the truncation
           응
               is performed at order `m`.
               **Resulting Dimensions:**
              If `A` is an `N×N` PhasorArray, then `BT(A, m)` is a
(2m+1)N \times (2m+1)N matrix.
           응
           응
               Syntax:
               r = BT(A, m)
                   Constructs a Block Toeplitz matrix of order `m`.
           양
           응
              Input Arguments:
               - ol (PhasorArray) : The input PhasorArray.
               - m (integer, optional) : The order of the Block Toeplitz
matrix.
           응
           응
               Output:
               - r ((2m+1)N \times (2m+1)N \text{ matrix}): The Block Toeplitz matrix.
           응
           응
              Example:
               % Generate a Block Toeplitz matrix from a PhasorArray
           응
               A = PhasorArray(rand(4,4,11));
               BT A = BT(A, 5);
              See also: TB, spTB.
           arguments
               01
               m = []
           end
           r=array2BToepliz(o1,2*m);
       end
        function r = TB(o1, m)
           %TB Construct a Toeplitz Block matrix from a PhasorArray.
               This function constructs a **Toeplitz Block (TB) matrix** of
order `m`, which is an alternative
               representation of `A` where **each block** of the matrix is
a Toeplitz structure.
               Specifically, TB(A, m) is given by:
                B = sum \{k=-2m\}^{2m} (A k \otimes I^k),
              where:
               - `I^k` is the **diagonal selector matrix**, which has ones
           응
on its `k`-th diagonal.
              - `⊗` denotes the **Kronecker product**.
               Unlike **Block Toeplitz (BT) **, where **blocks are repeated
along diagonals**, here **each block itself**
              follows a **Toeplitz structure**. Both `BT(A, m)` and `TB(A,
m) ` are equivalent in the **harmonic space**,
           응
              up to a basis transformation.
           용
               This matrix represents the **finite-dimensional truncation**
```

```
of an
               **infinite-dimensional Toeplitz Block operator** acting on
\ell^2, with truncation at order `m`.
                **Resulting Dimensions:**
               If `A` is an `N×N` PhasorArray, then `TB(A, m)` is a
(2m+1)N \times (2m+1)N matrix.
            응
                Syntax:
                r = TB(A, m)
            응
                    Constructs a Toeplitz Block matrix of order `m`.
            응
            응
               Input Arguments:
               - o1 (PhasorArray) : The input PhasorArray.
               - m (integer, optional) : The order of the Toeplitz Block
matrix.
            응
            용
               Output:
            응
               - r ((2m+1)N \times (2m+1)N \text{ matrix}): The Toeplitz Block matrix.
               Example:
            응
               % Generate a Toeplitz Block matrix from a PhasorArray
               A = PhasorArray(rand(4,4,11));
               TB A = TB(A, 5);
               See also: BT, spTB.
            arguments
                01
                m = []
            r=array2TBlocks(o1,2*m);
        end
        function r = spTB(o1, m)
            %SPTB Construct a sparse Toeplitz Block representation of A.
               This function generates a **sparse** Toeplitz Block
representation
               of order `m` for faster computations. The structure follows:
            응
                    B = sum \{k=-2m\}^{2m} (A k \otimes I^{k}),
               where:
               - `I^k` is the **diagonal selector matrix**, which has ones
on its `k`-th diagonal.
            % - `⊗` denotes the **Kronecker product**.
                - `A k` are the Fourier components (harmonics) of `A`.
               This represents the **truncation of an infinite-dimensional
            응
Toeplitz Block operator on \ell^{2**}.
               **Resulting Dimensions:**
               If `A` is an `N×N` PhasorArray, then `spTB(A, m)` is a
(2m+1)N \times (2m+1)N **sparse matrix**.
```

```
응
               Syntax:
                r = SPTB(A, m)
                    Computes a sparse Toeplitz Block representation of order
`m`.
            응
            응
                Input Arguments:
                - o1 (PhasorArray) : The input PhasorArray.
            응
                - m (integer, optional) : The order of the Toeplitz Block.
            응
                Output:
                - r ((2m+1)N \times (2m+1)N \text{ sparse matrix}): The sparse Toeplitz
Block matrix.
            응
               Example:
               % Generate a sparse Toeplitz Block matrix
            응
               A = PhasorArray(rand(4,4,11));
                spTB A = spTB(A, 5);
            응
                See also: TB, BT.
            arguments
                01
                m = []
            end
            r=sparray2TBlocks(o1,2*m);
        end
        function r = spBT(o1, m)
            %WIP%%%%%%%%%%%
            arguments
                01
                m = []
            end
            warning("NOT IMPLEMENTED YET")
                          r=sparray2BToeplitz(o1,2*m);
            r=sparse(array2BToepliz(o1,2*m));
        end
        function r = FvTB(o1, h)
            %FVTB Compute the Fourier representation of the vectorized form
of A(t).
            응
                This function first **vectorizes** the time-dependent matrix
            응
A(t) by stacking
                all its columns into a **single-column vector** a(t) =
vect(A(t)), and then
                applies the **TF TB** transformation to obtain its Fourier
            응
representation.
                **Alternative Computation:**
               - In MATLAB notation, vectorizing A(t) is equivalent to
**A{:}**.
                - Thus, this function is **equivalent to applying TF TB to
A{:}**, i.e.:
            응
                    FvTB(A, h) = TF TB(A\{:\}, h).
            응
                **Procedure:**
```

```
% 1. Compute the column-wise vectorization:
                    - If A(t) is an `N×M` matrix, then vect(A(t)) is a
column vector of size `NM \times 1`.
            This stacking is done **column by column** (following)
MATLAB's column-major order).
            % 2. Compute the Fourier series coefficients **up to order h**:
                   - Instead of applying `TF TB` directly to A(t), we apply
it to a(t) = vect(A(t)).
                   - The result is the Fourier representation of
`vect(A(t))`, which is a vectorized form
                   of TF TB(A, h).
               **Key Property:**
              If y(t) = A(t)x(t), then:
                   FvTB(y) = FvTB(A) \cdot FvTB(x),
              where `FvTB(A, h) ` is the **Fourier vectorized
representation ** of A.
                **Dimension of the Output:**
               - If `A` is an `N×M` matrix, then `FvTB(A, h)` is a
((2h+1)NM \times 1) vector.
            응
            응
               Syntax:
            응
               r = FVTB(A, h)
                   Computes the Fourier representation of `vect(A)` up to
order `h`.
               Input Arguments:
               - ol (PhasorArray) : The input PhasorArray representing A(t).
                - h (integer) : The highest harmonic order to retain in the
Fourier series.
               Output:
               - r ((2h+1)NM \times 1 \text{ vector}): The Fourier representation of
vect(A).
            응
            용
               Example:
               % Compute Fourier-vectorized representation of A(t)
               A = PhasorArray(rand(4,4,11));
            용
               FVTB A = FVTB(A, 5);
                \mbox{\%} Compute using TF TB on the vectorized form
               FvTB A alt = TF TB(A{:}, 5); % Equivalent computation
            9
               See also: TF TB, vect, trunc, pad.
            if size (o1, 3) > 2*h+1
                o1=trunc(o1,h);
            else
                o1=pad(o1,h-o1.h);
            end
            d1=reshape(o1,[],1,size(o1,3));
            r=pvalue(permute(d1,[3 1 2]));
            r=r(:);
```

end

```
function [HpJ, JHm, Hp, Hm] = TBHankel(o1, m)
            %TBHANKEL Compute Toeplitz Block Hankel matrices of order m.
                This function computes the **Toeplitz Block Hankel
matrices** (H) and the
                associated **J-Hankel matrices** (H J) of the given matrix
`A(t)` represented
               by the PhasorArray `o1`.
            응
               **Definition:**
                - A **Hankel matrix** is symmetric with entries reflected
across its antidiagonals.
               - A **J-Hankel matrix** is defined by pre- and post-
            용
multiplying the Hankel matrix
                  with the flipping operator \ \ J\ m\ \ , where \ \ \ J\ m\ \  flips the
diagonals of a matrix.
                **Key Property:**
                - For a periodic matrix function A(t), the Toeplitz Block
Hankel decomposition
            응
                  complements the Toeplitz Block transform (TB).
                **Truncation:**
                - The matrix is truncated to order `2m` harmonics before
            응
performing the computation.
            응
                **Output Details:**
               - `HpJ`: Positive J-Hankel matrix for `A(t)`.
                - `JHm`: Negative J-Hankel matrix for `A(t)`.
                - `Hp`: Positive Hankel block matrix for `A(t)`.
                - `Hm`: Negative Hankel block matrix for `A(t)`.
                **Dimensions:**
                - If `A` is an `N×N` matrix, then all outputs have
dimensions (2m+1)N \times (2m+1)N.
            응
                Syntax:
                [HpJ, JHm, Hp, Hm] = TBHANKEL(A, m)
                    Computes the Toeplitz Block Hankel and J-Hankel matrices
of order `m`.
            응
                Input Arguments:
                - ol (PhasorArray) : The input PhasorArray representing A(t).
                - m (integer) : The truncation order for the harmonics.
            응
            응
               Output:
                - HpJ ((2m+1)N × (2m+1)N matrix) : Positive J-Hankel matrix.
                - JHm ((2m+1)N \times (2m+1)N \text{ matrix}): Negative J-Hankel matrix.
               - Hp ((2m+1)N \times (2m+1)N \text{ matrix}): Positive Hankel block
matrix.
               - Hm ((2m+1)N × (2m+1)N matrix) : Negative Hankel block
matrix.
```

```
Example:
               % Compute Hankel matrices for a given A(t)
            % A = PhasorArray(rand(4,4,11));
               [HpJ, JHm, Hp, Hm] = TBHankel(A, 5);
               See also: spTBHANKEL, BTHANKEL.
            [HpJ, JHm, Hp, Hm] = Array2TBHankel (o1, 2*m);
        end
        function [HpJ, JHm, Hp, Hm] = spTBHankel(o1, m)
            %SPTBHANKEL Compute sparse Toeplitz Block Hankel matrices of
order m.
               This function computes the **sparse Toeplitz Block Hankel
matrices** (H) and
              their associated **J-Hankel matrices** (H J) for the given
PhasorArray `o1`.
               **Sparse Computation:**
              - The matrices are stored and computed in **sparse format**
for efficiency,
           응
                especially for high-dimensional problems or large
truncation orders.
               **Output Details:** (same as `TBHANKEL`)
              - `HpJ`: Positive sparse J-Hankel matrix.
            % - `JHm`: Negative sparse J-Hankel matrix.
               - `Hp`: Positive sparse Hankel block matrix.
               - `Hm`: Negative sparse Hankel block matrix.
            응
               **Syntax:**
               [HpJ, JHm, Hp, Hm] = SPTBHANKEL(A, m)
                    Computes the sparse Toeplitz Block Hankel and J-Hankel
matrices of order `m`.
               See also: TBHANKEL, BTHANKEL.
            [HpJ, JHm, Hp, Hm] = spArray2TBHankel(o1, 2*m);
        function [HpJ, JHm, Hp, Hm] = BTHankel(o1, m)
            %BTHANKEL Compute Block Toeplitz Hankel matrices of order m.
              This function computes the **Block Toeplitz Hankel
matrices** (H) and the
            % associated **J-Hankel matrices** (H J) for the given
PhasorArray `o1`.
                **Difference with TBHankel:**
              - While `TBHankel` corresponds to **Toeplitz Block (TB) **
structures, `BTHankel`
                corresponds to **Block Toeplitz (BT)** structures.
            응
              **Output Details:** (same as `TBHANKEL`)
              - `HpJ`: Positive J-Hankel matrix.
                - `JHm`: Negative J-Hankel matrix.
```

```
- `Hp`: Positive Hankel block matrix.
               - `Hm`: Negative Hankel block matrix.
               **Syntax:**
            응
            응
               [HpJ, JHm, Hp, Hm] = BTHANKEL(A, m)
                    Computes the Block Toeplitz Hankel and J-Hankel matrices
of order `m`.
                See also: SPBTHANKEL, TBHANKEL.
            [HpJ, JHm, Hp, Hm] = Array2BTHankel(o1, 2*m);
        end
        function [HpJ, JHm, Hp, Hm] = spBTHankel(o1, m)
            %SPBTHANKEL Compute sparse Block Toeplitz Hankel matrices of
order m.
              This function computes the **sparse Block Toeplitz Hankel
matrices** (H) and
              the associated **J-Hankel matrices** (H J) for the given
PhasorArray `o1`.
               **Sparse Computation:** (same as `spTBHANKEL`)
              - The matrices are stored and computed in **sparse format**
for efficiency.
               **Syntax:**
            응
            % [HpJ, JHm, Hp, Hm] = SPBTHANKEL(A, m)
                   Computes the sparse Block Toeplitz Hankel and J-Hankel
matrices of order `m`.
            % See also: BTHANKEL, TBHANKEL.
            [HpJ, JHm, Hp, Hm] = spArray2BTHankel(o1, 2*m);
        end
        function AB TB = TBmtimes(o1,o2,h)
            % TBMTIMES Compute the Toeplitz Block (TB) matrix of the product
A(t)B(t).
            응
               TBMTIMES(o1, o2, h) computes the Toeplitz Block matrix of
order `h`
               that represents the product of two periodic matrix functions
A(t) and B(t).
              Inputs:
                ol - (PhasorArray) The first input PhasorArray
representing A(t).
            % o2 - (PhasorArray) The second input PhasorArray
representing B(t).
                h - (integer) The truncation order for the harmonics.
            응
              Outputs:
                 AB TB - ((2h+1)N \times (2h+1)N \text{ matrix}) The Toeplitz Block
matrix for A(t)B(t).
              Behavior:
```

```
- Uses the Hankel matrices and TB representation of A(t)
and B(t).
                  - The resulting TB matrix satisfies:
                      T \text{ tb}(A * B) = T \text{ tb}(A) * T \text{ tb}(B) + H p(A) * H m(B) +
JH m(A) * JH p(B).
            응
               Output Dimensions:
                - If A and B are `N×N` matrices and the truncation order
is `h`, then:
                      - `AB TB` is a `(2h+1)N \times (2h+1)N` matrix.
            응
            응
               Example Usage:
                 % Compute TB matrix of the product A(t)B(t)
            응
            응
                  A = PhasorArray(rand(4,4,11));
            응
                 B = PhasorArray(rand(4, 4, 11));
            응
                  AB TB = TBmtimes(A, B, 5);
                See also: spTBmtimes, TB, TBHankel.
            [\sim, JHm, Hp, \sim] = TBHankel(o1,h);
            [HpJ2, \sim, \sim, Hm2] = TBHankel(o2, h);
            o1TB=o1.TB(h);
            o2TB=o2.TB(h);
            AB TB=o1TB*o2TB+Hp*Hm2+JHm*HpJ2;
        function AB TB = spTBmtimes(o1,o2,h)
            % SPTBMTIMES Compute the sparse Toeplitz Block (TB) matrix of
the product A(t)B(t).
            응
               SPTBMTIMES(01, 02, h) computes the sparse Toeplitz Block
matrix of order `h`
               that represents the product of two periodic matrix functions
A(t) and B(t).
            응
            % Inputs:
                 ol - (PhasorArray) The first input PhasorArray
representing A(t).
            응
                o2 - (PhasorArray) The second input PhasorArray
representing B(t).
            % h - (integer) The truncation order for the harmonics.
            응
               Outputs:
                AB TB - ((2h+1)N × (2h+1)N sparse matrix) Sparse TB matrix
for A(t)B(t).
            응
               Behavior:
                - Uses sparse versions of Hankel matrices and TB
representation of A(t) and B(t).
            % - The result satisfies:
                      T tb(A * B) = T tb(A) * T tb(B) + H p(A) * H m(B) +
JH m(A) * JH p(B).
            응
               Output Dimensions:
                - If A and B are `N×N` matrices and the truncation order
is `h`, then:
```

```
- `AB TB` is a `(2h+1)N × (2h+1)N` sparse matrix.
            응
            응
                Example Usage:
            응
                  % Compute sparse TB matrix of the product A(t)B(t)
            응
                  A = PhasorArray(rand(4,4,11));
            응
                  B = PhasorArray(rand(4,4,11));
            응
                  AB TB = spTBmtimes(A, B, 5);
                See also: TBmtimes, spTB, spTBHankel.
            [\sim, JHm, Hp, \sim] = spTBHankel(o1,h);
            [HpJ2, \sim, \sim, Hm2] = spTBHankel(o2, h);
            o1TB=o1.spTB(h);
            o2TB=o2.spTB(h);
            AB TB=o1TB*o2TB+Hp*Hm2+JHm*HpJ2;
        end
        function r= TF TB(o1,m)
            %TF TB Compute the Fourier representation of A in a form
compatible with TB(A, m).
                This function computes the **Fourier representation** of
            응
A(t) up to order `m`,
            용
               but instead of forming a **square Toeplitz Block matrix**
(TB(A, m)), it **stacks
               the phasors of each element of A into a structured
vectorized form**.
                **Definition:**
                - Let `a(t)` be a scalar function with Fourier coefficients
`(a k) \{k\in\mathbb{Z}\}`.
                - Then, TF TB(a, m) = [a (-m); a (-m+1); ... a 0; a 1; ...
a m]`.
                - If `A(t)` is an `N×M` matrix, its Fourier representation
is given by:
                    TF TB(A, m) = [TF TB(A 11), TF TB(A 12); TF TB(A 21),
TF TB(A 22)].
                **Key Property (TB Compatibility):**
                If \dot{y}(t) = A(t)x(t), then:
                    F TB(y) = TB(A, m) \cdot F TB(x),
                where `TB(A, m)` is the **Toeplitz Block representation** of
`A` (see `TB` method).
                **Dimension of the Output:**
                - If `A` is an `N×M` matrix, then `TF TB(A, m)` is a
((2m+1)N) \times M matrix.
            응
            응
               Syntax:
                r = TF TB(A, m)
                    Computes the Fourier representation of `A` up to order
m, compatible with TB(A, m).
            응
               Input Arguments:
                - ol (PhasorArray) : The input PhasorArray representing A(t).
```

```
% - m (integer, optional) : The highest harmonic order to
retain. Default: `A.h`.
            응
            응
               Output:
            % - r ((2m+1)N \times M \text{ matrix}): The Fourier representation of A
in a form compatible with TB(A, m).
            응
            응
                Example:
                % Compute Fourier representation of A in TB form
                A = PhasorArray(rand(4,4,11));
               TF TB A = TF TB (A, 5);
               See also: TF BT, TB.
            arguments
                01
                m = (size(o1, 3) - 1)/2;
            end
            nho=(size(o1,3)-1)/2;
            if nho<m</pre>
                toto=value(o1.pad(m-nho));
            elseif nho>m
                toto=pvalue(trunc(o1,m));
            else
                toto=pvalue(o1);
            end
            titi=permute(toto,[3 1 2]);
            r=reshape(titi,[],size(o1,2),1);
        end
        function r= TF BT(o1,m)
            %TF BT Compute the Fourier representation of A in a form
compatible with BT(A, m).
            응
                This function computes the **Fourier representation** of
A(t) up to order `m`,
               but instead of forming a **square Block Toeplitz matrix**
            응
(BT(A, m)), it **stacks
                the phasors of each element of A into a structured
vectorized form**.
                **Definition:**
               - Let `a(t)` be a scalar function with Fourier coefficients
`(a k) \{k\in\mathbb{Z}\}`.
                - Then, `TF BT(a, m) = [a (-m); a (-m+1); ... a 0; a 1; ...
a m]`.
                - If `A(t)` is an `N×M` matrix, its Fourier representation
is given by:
                    TF BT(A, m) = [TF BT(A 11), TF BT(A 12); TF BT(A 21),
TF BT(A 22)].
                **Key Property (BT Compatibility):**
            % If \dot{y}(t) = A(t)x(t), then:
                    F BT(y) = BT(A, m) \cdot F BT(x),
                where `BT(A, m)` is the **Block Toeplitz representation** of
`A` (see `BT` method).
```

```
응
                **Dimension of the Output:**
                - If `A` is an `N\timesM` matrix, then `TF BT(A, m)` is an `N \times
(2m+1)M matrix.
            응
                Syntax:
                r = TF BT(A, m)
                    Computes the Fourier representation of `A` up to order
m, compatible with BT(A, m).
            응
               Input Arguments:
               - ol (PhasorArray) : The input PhasorArray representing A(t).
                - m (integer, optional) : The highest harmonic order to
retain. Default: `A.h`.
            응
                Output:
            응
                - r (N \times (2m+1)M matrix) : The Fourier representation of A
in a form compatible with BT(A, m).
            응
            응
                Example:
            응
                % Compute Fourier representation of A in BT form
                A = PhasorArray(rand(4, 4, 11));
                TF BT A = TF BT (A, 5);
            응
               See also: TF TB, BT.
            arguments
                01
                m = (size(o1, 3) - 1)/2;
            end
            nho = (size(o1, 3) - 1)/2;
            if nho<m</pre>
                toto=padarray(pvalue(o1),[0 0 m-nho]);
            elseif nho>m
                toto=pvalue(trunc(o1,m));
            else
                toto=pvalue(o1);
            end
            titi=permute(toto,[1 3 2]);
            r=reshape(titi,[],size(o1,2),1);
        end
        function r= HmqNEig(o1,h,T,bandlimit)
            % HMQNEIG Compute the eigenvalues of TB(o1, h) - NTB(o1, h, T).
응
   HMQNEIG(o1, h, T, bandlimit) computes the eigenvalues of the difference
between
    the Toeplitz Block matrix `TB(o1, h)` and the non-periodic Toeplitz
approximation
응
    `NTB(o1, h, T)`, capturing spectral properties of `o1`.
응
용
    Inputs:
      01
                - (PhasorArray) The input PhasorArray representing a
periodic matrix function.
                - (integer, optional) The truncation order for the harmonics.
```

```
- Default: `2*o1.h` (twice the highest harmonic stored
in `o1`).
                - (double) The period of the PhasorArray.
      bandlimit - (char) Specifies band-limiting of eigenvalues:
응
응
                    - 'none' (default): No band-limiting.
                    - 'fundamental': Retains eigenvalues where |imag(r)| <
응
\pi / |T|.
응
응
    Outputs:
    r - (vector) The eigenvalues of the matrix `TB(o1, h) - NTB(o1, h, T)`.
응
응
   Behavior:
     - Computes eigenvalues of the difference matrix `Hmq = TB(o1, h) -
응
NTB(o1, h, T)`.
     - Band-limiting removes eigenvalues outside the specified range.
응
응
    Example Usage:
응
     % Compute eigenvalues with fundamental band-limiting
      A = PhasorArray(rand(4,4,11));
      r = HmqNEig(A, 10, 2, 'fundamental');
응
응
응
      % Compute eigenvalues without band-limiting
응
      r = HmqNEig(A, [], 1, 'none');
응
  See also: TB, NTB, eig.
            arguments
                01
                h = 20
                T=1
                bandlimit {mustBeMember(bandlimit,
{'none','fundamental'})}='none'
            end
            if isempty(h)
                h=2*o1.h;
            r1=o1.TB(h)-NTB(o1.value,h,T);
            r=eig(r1);
            switch bandlimit
                case 'fundamental'
                    r=r(abs(imag(r)) < pi/abs(T));
                case 'none'
                otherwise
                    r=r(abs(imag(r)) <=bandlimit);</pre>
            end
        end
        function r= HmgEig(o1,h)
            % HMQEIG Compute the eigenvalues of the Toeplitz Block (TB)
matrix of A(t).
                HMQEIG(01, h) computes the eigenvalues of the Toeplitz Block
(TB) matrix
                representation of the periodic matrix function `A(t)`,
capturing its spectral properties.
```

```
응
               Inputs:
                  o1 - (PhasorArray) The input PhasorArray representing a
periodic matrix function.
                h - (integer, optional) The truncation order for the
harmonics.
                        - Default: `ol.h` (the highest harmonic stored in
`o1`).
              Outputs:
              r - (vector) The eigenvalues of the Toeplitz Block
matrix `TB(o1, h)`.
               Behavior:
                - The matrix is truncated to order `h` to form `TB(o1, h)`.
                - If `h` is not specified, the function uses the highest
harmonic available in `o1`.
               Output Dimensions:
                - If `A(t)` is an `N×N` matrix and the truncation order is
`h`, then:
                      - `TB(o1, h)` is a `(2h+1)N × (2h+1)N` matrix.
            응
                      - `r` is a vector of eigenvalues of length `(2h+1)N`.
               Example Usage:
                 % Compute eigenvalues of TB(A, h) with a specific
truncation order
                 A = PhasorArray(rand(4,4,11));
            응
                 r = HmqEig(A, 10);
                 % Compute eigenvalues using the default harmonic order
                 r = HmqEiq(A);
              See also: TB, eig.
            arguments
               01
               h=o1.h
            end
            r1=o1.TB(h);
            r=eig(r1);
        end
        function Tout = stem(o1, varopt)
            % STEM Generate a stem plot for one or more `PhasorArray`
objects.
               STEM(o1, varopt) generates a stem plot to visualize the
phasors of one or more
               `PhasorArray` objects. Each input object is plotted with a
distinct marker style,
              and the function integrates seamlessly with existing
figures, axes, or tiled layouts.
```

```
% Inputs:
                             - (Repeating, `PhasorArray`) One or more
                01
`PhasorArray` objects to be plotted.
              Name-Value Pair Arguments:
                            - (char) Y-axis scale. Options:
           용
                                 - 'log' (default): Logarithmic scale.
                                 - 'linear': Linear scale.
                  'explosed' - (logical) Determines subplot arrangement:
                                 - true (default): Each matrix component is
plotted in its own subplot.
                                 - false: All components are combined in a
single plot.
                 'display' - (char) Specifies which part of the phasor to
display:
                                 - 'real': Real part.
                                 - 'imag': Imaginary part.
                                 - 'both': Both real and imaginary parts.
                                 - 'abs' (default): Magnitude of the
phasors.
                 'marker' - (cell array of chars) Marker styles for each
`PhasorArray`.
                                 - Default:
{"o","*","x","square","diamond","^","v",">","<"}.
                             - (char) Determines which harmonics are
               'side'
displayed:
                                 - 'both': Includes both positive and
negative harmonics.
                                 - 'oneSided' (default): Displays only non-
negative harmonics.
                 'parent' - (graphics handle) Parent figure or axes for
the plot.
                                 - Default: gcf (current figure).
              Outputs:
                             - (tiledlayout object) Handle to the tiled
           응
                 Tout
layout used for the plot.
              Behavior:
               - Multiple `PhasorArray` Objects: Each object in `o1` is
plotted separately using unique markers.
                  If more objects are provided than marker styles, markers
are cycled.
                - Real vs. Complex Phasors: Automatically adjusts `side`
to 'both' if any input contains complex values.
              Example Usage:
               % Plot absolute values of multiple PhasorArray objects
with linear scale
                A1 = PhasorArray(rand(3,3,11));
                 A2 = PhasorArray(rand(3,3,11));
                 Tout = stem(A1, A2, 'scale', 'linear', 'display', 'abs');
           용
                 % Plot real part of a PhasorArray using custom markers
```

```
A = PhasorArray(rand(4, 4, 11));
                  stem(A, 'display', 'real', 'marker', {"o", "^"});
            응
                  % Combine all components of multiple PhasorArray objects
in one plot
                  stem(A1, A2, 'explosed', false, 'side', 'both');
            응
                See also: stemPhasor, tiledlayout.
            arguments (Repeating)
                01
            end
            arguments
                varopt.scale {mustBeMember(varopt.scale,
{'log','linear'})}='log'
                varopt.explosed = true
                varopt.display {mustBeMember(varopt.display,
{'real', 'imag', 'both', 'abs'})} = 'abs'
                varopt.marker
={"o","*","x","square","diamond","^","v",">","<"};
                varopt.side {mustBeMember(varopt.side, {'both', 'oneSided'})}
= 'oneSided'
                varopt.parent = gcf
            end
            %check if all phasorArray in the cell o1 are real using cellfun
and isreal method
            if ~all(cellfun(@(x) isreal(x),o1))
                varopt.side='both';
            end
            if ~isa(varopt.marker, "cell")
                varopt.marker= { varopt.marker};
            end
            varhold=ishold;
stemPhasor(o1{1}, scale=varopt.scale, hold=varhold, explosed=varopt.explosed, dis
play=varopt.display, marker=varopt.marker{1}, side=varopt.side, parent=varopt.pa
rent);
            n=numel(o1);
            nmarker=numel(varopt.marker);
            for n iter=2:n
                oi=o1{n iter};
                ni=mod(n iter,nmarker);
                if mod(ni,nmarker) == 0
                    ni=nmarker;
                end
stemPhasor(oi,scale=varopt.scale,hold=true,explosed=varopt.explosed,marker=va
ropt.marker{ni}, display=varopt.display, parent=T);
            end
            hold off
            if varhold
                hold on
            end
```

```
if nargout>0
                Tout = T;
            end
        end
        function barsurf(o1,thres,hdel,varg)
            % BARSURF Generate a 3D bar surface plot of the phasors of a
`PhasorArray`.
               BARSURF(o1, thres, hdel, varg) produces a 3D bar surface
plot representing
               the phasors of `o1`. Phasors below a specified threshold are
            응
truncated, except
               for the first `hdel` harmonics that meet the condition,
            응
which are retained as a margin.
            응
            응
              Inputs:
                01
                         - (PhasorArray) The PhasorArray object to be
plotted.
                 thres - (double, optional) The relative threshold for
truncating phasors.
                             - Default: `1e-6`.
            응
                 hdel
                         - (integer, optional) The number of harmonics to
retain as a margin.
                             - Default: `3`.
                 varg - (struct, optional) Name-value pair arguments:
            응
            응
                             - 'scale' (char): Scale of the plot.
            응
                                 - 'log' (default): Logarithmic scale.
            응
                                 - 'linear': Linear scale.
            응
               Behavior:
                  - **Thresholding**: Phasors with absolute values below
`thres * maxPhasor`
            응
                   are truncated, except for the first `hdel` harmonics.
                  - **Logarithmic Scaling**: By default, the `log` scale is
used for better
                   visualization of magnitude variations.
                  - **Matrix Reshaping**: The function reshapes the phasors
into a suitable
                   format for `bar3` visualization.
            응
            응
            응
                Example Usage:
            용
                  % Generate a bar plot with default threshold and margin
            응
                 barsurf(o1);
            양
                  % Use a custom threshold and linear scaling
            응
                 barsurf(o1, 1e-6, 3, 'scale', 'linear');
                See also: ReduceArray, bar3.
            arguments
                thres=1e-6;
               hdel=3
                varg.scale {mustBeMember(varg.scale, {'log', 'linear'})} = 'log'
            end
```

```
if isa(o1.value, 'ndsdpvar')
                o1=value(value(o1));
                boolnan=isnan(o1);
                nnz (boolnan);
                if nnz(boolnan)>0
                    warning("Some sdpvar value are NaN")
                    o1(boolnan)=0;
                    o1=ReduceArray(o1);
                end
            end
            [nx, nz] = size(o1, [1 2]);
            nh = (size(o1, 3) - 1)/2;
[~, refM, hresM] = ReduceArray (o1, reduceMethod="relative", reduceThreshold=thres, e
xcludeOPhasor=false);
            %refM contient en val absolue le plus grand phasor de chaque
            %composante de o1.
            minM signif=min(abs(refM),[],'all')*thres;
            hdel=min(nh-hresM,hdel);
            epsM=10^{(floor(log10(minM signif)))*10^-(1.5)};
            reshM=abs(reshape(ReduceArray(o1,hresM+hdel),nx*nz,[]));
            barsurf(reshM(:,((end+1)/2):end).',epsM,"yticklabel",
(0:hresM+hdel)','scale',varg.scale)
                 barsurf(reshM(:,hres+1:end),min(ref,
[], 'all') *thres, "xticklabel", (0:hres)', 'scale', 'log')
            xlabel("States")
            ylabel("Harmonics")
            title('Phasor of M')
            xlim([0 nx*nz+1])
        end
        function [r,t]=plot(o1,T,t,arg)
            % PLOT Evaluate and plot a T-periodic `PhasorArray`.
               PLOT(o1, T, t, arg) computes and plots the time-domain
representation of a
                `PhasorArray`, assuming it is `T`-periodic.
               This method evaluates the time-domain representation of a
`PhasorArray` and
              plots it over a specified period `T`. It provides flexible
options for
               visualization, including real/imaginary decomposition,
subplot arrangements,
            응
                and axis linking.
            응
                Inputs:
                  ol - (PhasorArray) The `PhasorArray` object to be
evaluated and plotted.
            응
                 T - (double, optional) Period of the `PhasorArray`.
                           - Default: `1`.
            응
                 t - (vector or scalar, optional) Time instants for
```

```
evaluation.
                          - If `t = []`: A default time grid is generated
as 0:dt:T-dt, where dt = T/(20*h), with h the highest harmonic.
                         - If `t = [tmin tmax]`: Uses `t = tmin:dt:tmax`
with `dt` computed as above.
                          - If `t` is a vector: Evaluates `A(t)` at the
specified values.
                          - If `t` is a scalar: Evaluates `A(t)` at that
single instant.
            % arg - (struct, optional) Name-value pair arguments:
                          - 'plot' (logical): Display the plot (default:
`true`).
                          - 'explosed' (logical): Plot each matrix
component in a separate subplot (default: `true`).
                          - 'hold' (logical): Hold the current plot
(default: `false`).
                          - 'DispImag' (logical): Display the imaginary
part of the matrix (default: `false`).
                          - 'DispReal' (logical): Display the real part of
the matrix (default: `true`).
                          - 'ZeroCentered' (logical): Center the Y-axis
around zero (default: `false`).
           응
                          - 'title' (string): Custom title for the figure
(default: `[]`).
                          - 'linetype' (string): Line style for the plot
(default: `'-'`).
                          - 'GlobalYLim' (logical): Apply the same Y-limits
across subplots (default: `false`).
                          - 'linkaxes' (string): Link axes of subplots
(`'x'`, `'y'`, `'xy'`, etc.) (default: `'x'`).
                          - 'forceReal' (logical): Assume `o1` is real-
valued and simplify computation (default: `false`).
           % Outputs:
                r - (m \times n \times length(t) array) Evaluated time-domain
representation of `o1`.
           % t - (vector) Time instants at which `o1` is evaluated.
              Behavior:
               - **Evaluation**: Computes `A(t)` using
           응
`PhasorArray2time`, which performs an inverse Fourier summation.
           % - **Automatic Time Grid**: If `t = []`, it generates a
default time grid based on `T` and harmonics.
           % - **Real-Valued Constraints**: If `o1` is real-valued,
`forceReal` is automatically set to `true`.
           % - **Plot Customization**:
                    - Supports separate subplots for each matrix element
(`explosed = true`).
                     - Can plot only the real or imaginary part
(`DispReal`, `DispImag`).
                    - Allows axis linking and uniform Y-axis scaling
across subplots.
            % Example Usage:
```

```
% Evaluate and plot a PhasorArray over one period
                  A = PhasorArray(rand(3,3,11));
                  plot(A, 2*pi, []);
            응
            응
                  % Evaluate A(t) on a custom time range
                  t = linspace(0, 2*pi, 100);
            응
                  plot(A, 2*pi, t, 'plot', true, 'explosed', false);
                  % Plot only the imaginary part with a different linestyle
                  plot(A, 2*pi, [], 'DispImag', true, 'DispReal', false,
'linetype', '--');
                See also: PhasorArray2time.
            arguments
                01
                T=1
                t=[]
                arg.plot logical =true
                arg.explosed logical =true
                arg.hold logical =false
                arg.DispImag logical =false
                arg.DispReal logical =true
                arg.ZeroCentered logical =false
                arg.title=[]
                arg.linetype='-'
                arg.GlobalYLim logical =false
                arg.linkaxes='x'
                arg.forceReal = false
            end
            if ishold
                arg.hold = true;
            end
            if isreal(o1)
                arg.forceReal = true;
            end
            [rr,tt]=PhasorArray2time(o1,T,t,plot=arg.plot,
DispImag=arg.DispImag, ...
DispReal=arg.DispReal,explosed=arg.explosed,hold=arg.hold,ZeroCentered=arg.Ze
roCentered, ...
title=arg.title,linetype=arg.linetype,GlobalYLim=arg.GlobalYLim,linkaxes=arg.
linkaxes, forceReal=arg.forceReal);
            if nargout>0
                r=rr;
            end
            if nargout>1
                t=tt;
            end
        end
        function r=plot3D(o1,T,t,arg)
            %PLOT3D Produce a 3D plot where x-axis is the real part, y-axis
is the imaginary part, and z-axis is time
```

```
응
                r = PLOT3D(o1, T, t, arg) produces a 3D plot where the x-
axis is the real part of A(t),
                the y-axis is the imaginary part of A(t), and the z-axis is
time.
                For each element of A, a subplot is created.
            응
               Input Arguments:
                o1 - The PhasorArray object to be plotted.
                T - The period of the PhasorArray. Default is 1.
                t - Time instants at which to evaluate the PhasorArray. Can
be:
                    - [] (empty), then t takes the value [0 T].
                    - [tmin tmax], then t=tmin:dt:tmax with automatic
discretization.
            응
                    - A vector on which A(t) is evaluated.
                    - A single scalar, then t takes the value [0 t].
                arg - (Optional) Name-value pair arguments:
                    'ZeroCentered' - Logical flag to normalize x and y axes
around zero. Default is false.
                    'title' - String to display a custom title to the
figure. Default is [].
                    'GlobalYLim' - Logical flag to enforce same Y and X
limits on the axes. Default is false.
                    'linkaxes' - String to link zoom on the x, y, z axes.
Default is 'x'.
               Output Arguments:
               r - The evaluated PhasorArray at the specified time instants.
               Example:
                r = plot3D(o1, 2*pi, []);
                r = plot3D(o1, 2*pi, [0 2*pi], 'ZeroCentered', true,
'title', '3D Plot');
                See also: PhasorArray2time
            arguments
                01
                T=1
                t=[0 T]
                arg.ZeroCentered=false
                arg.title=[]
                arg.GlobalYLim=false
                arg.linkaxes='x'
            end
            if numel(t) ==1
                t=sort([0 t]);
            rr=PhasorArray2time(o1, T, t, plot=true, plot3D=true,
DispImag=false,
DispReal=true, explosed=true, ZeroCentered=arg.ZeroCentered, title=arg.title,Glo
balYLim=arg.GlobalYLim,linkaxes=arg.linkaxes);
            if nargout
```

```
r=rr;
            end
        end
        function [r,t]=sim(o1,T,t,arg)
            % SIM Evaluate the time-domain representation of a `PhasorArray`.
               SIM(o1, T, t, arg) computes the time-domain representation
of a `PhasorArray`,
               assuming it is `T`-periodic.
            응
               SIM is a **convenience function** for evaluating `A(t)`,
leveraging `PhasorArray2time`.
           응
              It includes an option to enforce real-valued computation
when A(t) is known to be real.
            응
              Inputs:
            응
                o1 - (PhasorArray) The `PhasorArray` object representing
a periodic matrix.
                    - (double, optional) The period of `A(t)`.
                          - Default: `1`.
                t - (vector or scalar, optional) Time instants for
evaluation.
                           - If `t = []`: Uses a default time grid `0:dt:T-
dt', where 'dt = T/(20*h)', and 'h' is the highest harmonic.
                           - If `t = [tmin tmax]`: Uses `t = tmin:dt:tmax`
with automatic step size.
                          - If `t` is a vector: Evaluates `A(t)` at the
           9
specified values.
                          - If `t` is a scalar: Evaluates `A(t)` at a
single instant.
                  arg - (struct, optional) Name-value pair arguments:
                          - 'isRealValued' (logical): Enforce real-valued
computation (default: `false`).
            응
              Outputs:
                r - (m \times n \times length(t) array) Evaluated time-domain
representation of `o1`.
               t - (vector) Time instants at which `o1` was evaluated.
            응
              Behavior:
                - If `o1` is **real-valued**, `isRealValued` is
automatically set to `true`.
                 - Calls `PhasorArray2time` with a structured argument list
for evaluation.
            응
            응
               Example Usage:
                 % Evaluate A(t) over one period
                 A = PhasorArray(rand(3,3,11));
           용
                 [r, t] = sim(A, 2*pi, []);
                  % Force real-valued computation
            응
                 [r, t] = sim(A, 2*pi, [], 'isRealValued', true);
```

```
See also: PhasorArray2time, plot.
            arguments
                01
                T=1
                t=[]
                arg.isRealValued = false;
            end
            if isreal(o1)
                arg.isRealValued = true;
            end
            argo=struct;
            argo.plot=false;
            argo.explosed=true;
            argo.hold=false;
            argo.DispImag=false;
            argo.DispReal=true;
            argo.ZeroCentered=false;
            argo.title=[];
            argo.forceReal = arg.isRealValued;
            C=namedargs2cell(argo);
            [r,t]=PhasorArray2time(o1,T,t,C{:});
        function r=evalTime(o1,T,t)
            % EVALTIME Evaluate a `PhasorArray` in the time domain for a
given period `T`.
                EVALTIME(o1, T, t) computes the time-domain representation
of a `PhasorArray`,
            응
               assuming it is `T`-periodic. The function returns `A(t)`,
evaluated at specified
               time instants `t`, without generating any plots.
            응
                Inputs:
                  ol - (PhasorArray) The `PhasorArray` object to be
evaluated.
                      - (double, optional) The period of the `PhasorArray`.
                  Т
                            - Default: `1`.
                      - (vector or scalar, optional) Time instants for
                  t
evaluation.
                           - If `t = []`: Uses a default grid `0:dt:T-dt`,
with \dot{d}t = T/(20*h), where \dot{h} is the highest harmonic.
                           - If `t = [tmin tmax]`: Uses `t = tmin:dt:tmax`
with automatic step size.
                           - If `t` is a vector: Evaluates `A(t)` at the
specified values.
                           - If `t` is a scalar: Evaluates `A(t)` at a
single instant.
            응
                Outputs:
                  r - (m \times n \times length(t) array) Evaluated time-domain
```

응

```
representation of `o1`.
                Behavior:
                  - Uses `plot(o1, T, t, ...)` internally but **without
plotting** (`'plot', false`).
                  - Automatically generates a time grid if `t = []`.
            응
                  - Designed for numerical evaluation without visualization.
            응
            응
                Example Usage:
                  % Evaluate A(t) over one period
            응
                  A = PhasorArray(rand(3,3,11));
            응
                 r = evalTime(A, 2*pi, []);
            응
            응
                  % Evaluate A(t) on a custom time range
                  t = linspace(0, 2*pi, 100);
            응
                  r = evalTime(A, 2*pi, t);
                See also: plot, PhasorArray2time.
            arguments
                01
                T=1
                t=[]
            end
            argo=struct;
            argo.plot=false;
            argo.explosed=true;
            argo.hold=false;
            argo.DispImag=false;
            argo.DispReal=true;
            argo.ZeroCentered=false;
            argo.title=[];
            C=namedargs2cell(argo);
            r=plot(o1, T, t, C{:});
        end
        function [y,t,dy]=initial(o1,x0,T,tfinal,varg)
            % INITIAL Compute the system response to an initial state in a
periodic state-space model.
                INITIAL(o1, x0, T, tfinal, varg) simulates the system
response for:
                    dx/dt = A(t)x
               where `A(t)` is a `T`-periodic state-space matrix. This
            응
function calls `lsim` to compute the response.
            응
                Syntax:
                  [y, t] = INITIAL(A, x0, T, tfinal)
                  [y, t, dy] = INITIAL(A, x0, T, tfinal, Name, Value)
            응
            응
                Inputs:
                  01
                         - (PhasorArray) The time-varying system matrix
`A(t)`, stored as a **3D phasor array**.
            용
                  x0
                         - (vector, optional) Initial state `x(0)`.
            응
                             - Default: `ones(size(o1,1),1)`.
                         - (double, optional) Period of the system. Default:
```

```
`1`.
               tfinal - (scalar or vector, optional) Final simulation
time.
                            - Default: `10*T`.
           응
           응
                            - If scalar: Simulates from `t=0` to `t=tfinal`.
                            - If `[tmin tmax]`: Simulates from `tmin` to
           응
`tmax`.
                            - If vector: Uses provided time grid.
           용
              Name-Value Pair Arguments:
                 'opts'
           응
                           - (struct) Options for the ODE solver
(`odeset`). Default: `[]`.
                                   - (logical) Plot the state trajectory.
           응
                 'plot'
Default: `true`.
                                - (char) ODE solver method. Default:
           응
                 'solver'
`'adaptative'`. Options:
                                         - `'adaptative'`, `'forward-
euler'`, `'RK4'`
          응
                'FSprecpow'
                                - (integer) Power of 2 for frequency
sampling. Default: `8`.
                'checkReal'
                                 - (logical) Force real-valued output
(only if system is **guaranteed** real). Default: `false`.
          % 'isRealValued' - (logical) Force real-valued
computation. Default: `false`.
              Outputs:
                y - (matrix) State trajectory of the system (`size(y,1) =
size(o1,1)`).
                t - (vector) Time instants at which `y(t)` is evaluated.
                dy - (matrix, optional) Derivative of the state trajectory
(only if `nargout > 2`).
             Behavior:
                - This function is a **wrapper for `lsim`**, automatically
setting U(t) = 0.
           % - If `isRealValued` is **not provided**, it is
automatically detected from `o1`.
                - If `tfinal = 0`, the function **automatically sets**
`tfinal = 10*T`.
           응
           응
              Example Usage:
                % Simulate free response over one period
                A = PhasorArray(rand(3,3,11));
           응
           응
                [y, t] = initial(A, ones(3,1), 2*pi, 2*pi);
                 % Simulate with a fixed-step RK4 method
                [y, t] = initial(A, ones(3,1), 2*pi, 2*pi, 'solver',
'RK4');
               See also: LSIM, HMQ SIM.
           arguments
               x0 = ones(size(o1, 1), 1)
               T=1
```

```
varq.opts=[] %odeset option
                varg.plot=true
                varg.solver='adaptative'
                varg.FSprecpow=8
                varg.checkReal=0
                varg.isRealValued logical = false
            end
            if isreal(o1)
                varq.isRealValued = true;
            end
            vvarg = namedargs2cell(varg);
            if nargout == 3
                [y,t,dy] = lsim(o1,tfinal,x0,T,vvarg{:});
            else
                [y,t]=lsim(o1,tfinal,x0,T,vvarg\{:\});
            end
        function [y,t,dy]=lsim(o1,tfinal,x0,T,Uph,varg)
            % LSIM Simulate the response of a time-periodic linear system.
               LSIM(o1, tfinal, x0, T, Uph, varg) simulates the system:
                    dx/dt = A(t)x + U(t)
                where A(t) and U(t) are T-periodic matrices,
represented as phasor arrays.
                The function supports **adaptive and fixed-step solvers**,
allows **real-valued computation**, and
                provides **ODE solver customization** via `odeset` options.
            응
            응
                Syntax:
            용
                  [y, t] = LSIM(A, tfinal, x0, T)
                  [y, t, dy] = LSIM(A, tfinal, x0, T, U, Name, Value)
            응
            응
            응
                Inputs:
                  01
                         - (PhasorArray) The periodic system matrix `A(t)`,
stored as a **3D phasor array**.
                  tfinal - (scalar or vector, optional) Final simulation
time.
                             - Default: `10*T`
                             - If scalar: Simulates from `t=0` to `t=tfinal`.
            응
            응
                             - If `[tmin tmax]`: Simulates from `tmin` to
`tmax`.
                             - If vector: Uses provided time grid.
            응
                         - (vector, optional) Initial condition `x(0)`.
                  x0
                             - Default: `ones(size(o1,1),1)`.
                         - (double, optional) The period of `A(t)`. Default:
`1`.
                         - (PhasorArray or matrix, optional) The time-
                  Uph
varying input matrix `U(t)`.
            응
                             - Default: `[]` (zero input).
            응
                Name-Value Pair Arguments:
```

tfinal=0

```
% 'opts'
                                    - (struct) Options for the ODE solver
(`odeset`). Default: `[]`.
                                    - (logical) Plot the state trajectory.
                 'plot'
           응
Default: `true`.
                 'solver'
           응
                                   - (char) ODE solver method. Default:
`'adaptative'`. Options:
                                          - `'adaptative'`, `'forward-
           응
euler'`, `'RK4'`
           % 'FSprecpow' - (integer) Power of 2 for frequency
sampling. Default: `8`.
                 'checkReal'
                                   - (logical) Force real-valued output
           용
(only if system is **guaranteed** real). Default: `false`.
          % 'isRealValued' - (logical) Force real-valued
computation. Default: `false`.
           용
           % Outputs:
               y - (matrix) State trajectory of the system (`size(y,1) =
size(o1,1)`).
                t - (vector) Time instants at which `v(t)` is evaluated.
               dy - (matrix, optional) Derivative of the state trajectory
           응
(only if `nargout > 2`).
           용
              Behavior:
               - If `isRealValued` is **not provided**, it is
automatically detected from `o1` and `Uph`.
           % - **Default solver:** `ode15s` (adaptive).
                - If `tfinal = 0`, the function **automatically sets**
`tfinal = 10*T`.
                - If `Uph` is empty, the system is simulated as
           응
**homogeneous** (^dx/dt = A(t)x).
           용
               - The highest resolved harmonic `h` determines the
**integration step size**.
           용
           응
              Example Usage:
                % Simulate system response over one period
           응
           응
                A = PhasorArray(rand(3,3,11));
           응
                 [y, t] = lsim(A, 2*pi, ones(3,1), 2*pi);
           양
           응
                 % Simulate with a time-varying input U(t)
           응
                 U = PhasorArray(rand(3,1,11));
                 [y, t] = lsim(A, 2*pi, ones(3,1), 2*pi, U);
           응
           응
                % Use a fixed-step RK4 method
           응
                 [y, t] = lsim(A, 2*pi, ones(3,1), 2*pi, [], 'solver',
'RK4');
              See also: HMQ SIM, ode15s.
           arguments
               ^{\circ}1
               tfinal=0
               x0 = ones(size(o1, 1), 1)
               T=1
               Uph=[]
               varg.opts=[] %odeset option
```

```
varg.solver='adaptative'
                varg.FSprecpow=8
                varg.checkReal=0
                varg.isRealValued logical = false
            end
            if isempty(x0)
                x0 = ones(size(o1, 1), 1);
            end
            if isreal(o1)
                varg.isRealValued = true;
            end
            C=namedargs2cell(varg);
            C\{1\}="odeOpts";
            %asking derivative trigger more computation from hmq sim,
            %procede with care
            if nargout>2
                [y,t,dy] = hmq sim(o1,tfinal,x0,T,Uph,C{:});
            else %sinon
                [y,t]
                       = hmq sim(o1, tfinal, x0, T, Uph, C\{:\});
            end
        end
        function r=evalp(o1, angle, arg)
            % EVALP Evaluate a periodic matrix at a given phase angle.
                EVALP(01, angle, arg) computes the value of a periodic
matrix `A` at a
            응
               specified phase angle `0` instead of time. This function is
useful for
                evaluating periodic signals in the frequency domain.
            응
            응
            응
               Syntax:
                 r = EVALP(A, angle)
            응
                 r = EVALP(A, angle, Name, Value)
            응
            응
               Inputs:
                       - (PhasorArray) The periodic matrix represented as a
                 01
**phasor array**.
                  angle - (double, vector) The phase angle(s) at which to
            응
evaluate A(\theta).
               Name-Value Pair Arguments:
            응
                  'forceReal' - (logical) If `true`, forces real-valued
output. Default: `false`.
            응
                 'checkReal'
                                 - (logical) If `true`, checks whether the
result is real-valued. Default: `false`.
            %
                 'checkRealTol' - (double) Tolerance for checking real-
valued results. Default: `1e-8`.
              Outputs:
```

varg.plot=true

```
r - (matrix) Evaluated periodic matrix at the given
**phase angle** \theta.
            응
              Behavior:
                - This function evaluates `A(\theta)`, where `\theta` represents the
**phase angle** of the periodic function.
            % - The function internally calls `PhasorArray2time` with `T
= 2\pi, assuming **one full cycle** of the periodic function.
              - If `o1` is detected as real-valued, `forceReal` is
automatically set to `true`.
            응
               Example Usage:
                % Evaluate a phasor array at a phase angle of \pi/4
            응
                 result = evalp(A, pi/4);
            응
                % Force real-valued output
                 result = evalp(A, pi/4, 'forceReal', true);
            용
                See also: PHASORARRAY2TIME, EVALTIME.
            arguments
                01
                angle
                arg.forceReal logical
                                        = false
                                        = false
                arg.checkReal logical
                arg.checkRealTol
                                         = 1e-8
            end
            if isreal(o1)
                arg.forceReal = true;
            end
            varg = namedargs2cell(arg);
            %evaluate periodic matrix A for an angle argument instead of
            %time
            r=PhasorArray2time(o1,2*pi,angle,varg{:});
        end
        function r = mreal(01)
            %MREAL Compute the real part of a PhasorArray in the time domain.
                r = MREAL(01) returns the real part of the PhasorArray o1 in
the time domain.
                The real part R(t) is such that A(t) = R(t) + i*I(t), where
I(t) is the imaginary part.
               Both R(t) and I(t) are real-valued.
            응
            응
            응
                INPUT:
            응
                    ol - The PhasorArray object.
            응
            응
               OUTPUT:
            응
                  r - The real part of the PhasorArray in the time domain.
            응
            응
               Example:
                    A = PhasorArray.random(3, 3, 5);
```

```
% R = mreal(A);
                See also: mimag, mconj
            dval=pvalue(o1);
            r = real(dval + flip(dval, 3))*(1/2) + 1i * imag(dval -
flip(dval, 3)) *(1/2);
           r = PhasorArray(r);
        end
        function r = mimag(o1)
            %MIMAG Compute the imaginary part of a PhasorArray in the time
domain.
                r = MIMAG(o1) returns the imaginary part of the PhasorArray
ol in the time domain.
                The imaginary part I(t) is such that A(t) = R(t) + i*I(t),
where R(t) is the real part.
            응
              Both R(t) and I(t) are real-valued.
            응
            응
               INPUT:
                   ol - The PhasorArray object.
            응
            응
            응
               OUTPUT:
            응
                   r - The imaginary part of the PhasorArray in the time
domain.
            응
               Example:
            응
                   A = PhasorArray.random(3, 3, 5);
            응
                   I = mimag(A);
            응
            % See also: mreal, mconj
           dval=pvalue(o1);
            r = real(dval - flip(dval, 3))/2/1i + 1i * imag(dval +
flip(dval,3))/2/1i;
           r = PhasorArray(r);
        end
        function r = mconj(o1)
            %MCONJ Compute the complex conjugate of a PhasorArray in the
time domain.
              r = MCONJ(o1) returns the complex conjugate of the
PhasorArray ol in the time domain.
              The complex conjugate is computed as R(t) - i*I(t), where
R(t) is the real part and I(t) is the imaginary part.
           응
            응
                INPUT:
                   o1 - The PhasorArray object.
               OUTPUT:
            응
                    r - The complex conjugate of the PhasorArray in the
time domain.
            응
               Example:
                   A = PhasorArray.random(3, 3, 5);
            응
                   C = mconj(A);
```

```
응
        See also: mreal, mimag
    r = mreal(o1) - 1i * mimag(o1);
end
function o1=confirm reality(o1)
    % depreciated
    응
    % see also MREAL.
    pos part=o1{:,:,1:o1.h};
    neg part=o1{:,:,-1:-1:(-o1.h)};
    z part=o1.phas(0);
    o1{:,:,1:o1.h}=(pos part+conj(neg part))/2;
    o1{:,:,0}=real(z part);
    o1{:,:,-1:-1:(-o1.h)}=conj(pos part+conj(neg part))/2;
end
function r = pageconj(o1)
    %phasor conjugate
    r=(conj(pvalue(o1)));
end
function r = pagereal(o1)
    %phasor real part
    r=(real(pvalue(o1)));
end
function r = pageimag(o1)
    %phasor imag part
    r=(imag(pvalue(o1)));
function r = pageabs(o1)
    %phasor absolute value
    r=(abs(pvalue(o1)));
end
function r = isvector(o1)
    % ISVECTOR True if A(t) is a vector (row or column).
    r=isvector(o1.phas(0));
end
function r = isscalar(o1)
    % ISSCALAR True if A(t) is a scalar.
    r=isscalar(o1.phas(0));
end
function r = isrow(o1)
    % ISROW True if A(t) is a row vector.
    r=isrow(o1.phas(0));
end
function r = iscolumn(o1)
    % ISCOLUMN True if A(t) is a column vector.
    r=iscolumn(o1.phas(0));
end
function r = ismatrix(o1)
```

```
r=ismatrix(o1.phas(0));
        end
        function r = isnumeric(o1)
            % ISNUMERIC True if A(t) contains numeric values.
            r=isnumeric(o1.value);
        end
        function r = islogical(o1)
            % ISLOGICAL True if A(t) contains logical values.
            r=islogical(o1.value);
        end
        function r = isempty(o1)
            % ISEMPTY True if A(t) is empty.
            r=isempty(o1.value);
        function r = issquare(o1)
            % ISSQUARE True if A(t) is a square matrix.
            r = (size(o1,1) == size(o1,2));
        end
        function [r,R] = issymmetric(o1,arg)
            % ISSYMMETRIC Check if the PhasorArray object is symmetric.
                [r, R] = ISSYMMETRIC(o1, arg) checks if the PhasorArray
object o1
            응
               is symmetric based on the specified arguments.
            응
            응
               Input arguments:
                    o1 - The PhasorArray object to be checked.
                    arg - A structure with the following fields:
                        skewOption - A string specifying the type of
symmetry to check.
                                      It can be either 'nonskew' or 'skew'.
            응
Default is 'nonskew'.
                       tol - A tolerance value for the symmetry check.
Default is 0.
            % Output arguments:
                    r - A logical value indicating if the entire PhasorArray
object is symmetric.
            용
                    R - A logical array indicating the symmetry of each
slice of the PhasorArray object.
            응
               Example:
                  [r, R] = issymmetric(o1, 'skewOption', 'nonskew', 'tol',
1e-6);
               See also: arrayfun, norm, nnz
            arguments
                arg.skewOption {mustBeMember(arg.skewOption,
{'nonskew','skew'})} = 'nonskew'
                arg.tol = 0
            end
```

% ISMATRIX True if A(t) is a matrix.

```
if arg.tol == 0
                 R = arrayfun(@(ii) issymmetric(o1(:,:,ii),arg.skewOption),1:
(2*o1.h+1));
             else
                 tol=abs(arg.tol);
                 switch arg.skewOption
                     case 'nonskew'
                         R = \operatorname{arrayfun}(@(ii) \operatorname{norm}(o1(:,:,ii) -
o1(:,:,ii).',"inf")/norm(o1(:,:,ii),"inf")<to1,1:(2*o1.h+1));
                     otherwise
                         R = \operatorname{arrayfun}(@(ii) \operatorname{norm}(o1(:,:,ii))
+o1(:,:,ii).',"inf")/norm(o1(:,:,ii),"inf")<tol,1:(2*o1.h+1));
            end
             r = ((nnz(R)) == (2*o1.h+1));
        end
        function [r,R] = ishermitian(o1,arg)
             % ISHERMITIAN Check if the PhasorArray object is Hermitian.
                 [r, R] = ISHERMITIAN(o1, arg) checks if the PhasorArray
object o1
                is Hermitian. The function returns a logical scalar r
indicating
                 if all slices of the PhasorArray are Hermitian, and a
logical array R
             응
                 indicating if each individual slice is Hermitian.
            응
            응
                 Input arguments:
             응
                     o1 - PhasorArray object to be checked.
             응
                     arg - Structure with the following fields:
                         skewOption - (optional) Specifies whether to check
for
                                        'nonskew' (default) or 'skew'
Hermitian.
                                        Must be one of {'nonskew', 'skew'}.
                         tol - (optional) Tolerance for numerical comparison.
Default is 0.
                 Output arguments:
                    r - Logical scalar indicating if all slices of the
PhasorArray
                         are Hermitian.
                     R - Logical array indicating if each individual slice of
the
                         PhasorArray is Hermitian.
            arguments
                 01
                 arg.skewOption {mustBeMember(arg.skewOption,
{'nonskew', 'skew'})} = 'nonskew'
                 arg.tol = 0
            if arg.tol == 0
                 R = arrayfun(@(ii) ishermitian(o1(:,:,ii),arg.skewOption),1:
```

```
(2*o1.h+1));
            else
                tol=abs(arg.tol);
                switch arg.skewOption
                     case 'nonskew'
                         R = \operatorname{arrayfun}(@(ii) \operatorname{norm}(o1(:,:,ii) -
o1(:,:,ii)',"inf")/norm(o1(:,:,ii),"inf")<tol,1:(2*o1.h+1));
                     otherwise
                         R = \operatorname{arrayfun}(@(ii) \operatorname{norm}(o1(:,:,ii))
+o1(:,:,ii)',"inf")/norm(o1(:,:,ii),"inf")<tol,1:(2*o1.h+1));
            end
            r = ((nnz(R)) == (2*o1.h+1));
        end
        function [r,R] = isreal(o1,tol)
            % ISREAL Check if the imaginary part of the time realization of
            % PhasorArray is negligible compared to a given tolerance.
            응
            % Syntax:
                [r, R] = isreal(o1)
                [r, R] = isreal(o1, to1)
            % Description:
                This function checks if the imaginary part of the time
realization
                of a PhasorArray object is negligible compared to a
specified tolerance.
                It returns a boolean value 'r' indicating whether the
imaginary part
               is negligible, and a logical array 'R' indicating the same
for each
            용
               element of the PhasorArray.
            응
            % Inputs:
               o1 - PhasorArray object to be checked.
               tol - (Optional) Tolerance value for comparison. If not
provided,
            응
                       the function uses the default machine epsilon.
            응
            % Outputs:
               r - Boolean value (true/false) indicating if the imaginary
part
                       is negligible for the entire PhasorArray.
                R - Logical array indicating if the imaginary part is
negligible
                      for each element of the PhasorArray.
            % Example:
            % [r, R] = isreal(o1)
                [r, R] = isreal(o1, 1e-6)
            % See also:
            % mreal, ismembertol, abs
            arguments
```

```
01
                tol=[]
            end
            o1 r = mreal(o1);
            %if o1.value is not sym, sdpvar nor ndsdpvar
            if ~(isa(o1.value, "sym") || isa(o1.value, "ndsdpvar") ||
isa(o1.value, "sdpvar"))
                if isempty(tol)
                    r1 = ismembertol(real(o1.value), real(o1 r.value));
                    r2 = ismembertol(imag(o1.value),imag(o1 r.value));
                else
                    r1 = ismembertol(real(o1.value), real(o1 r.value), tol);
                    r2 = ismembertol(imag(o1.value),imag(o1 r.value),tol);
                end
            else %in this case ismembertol doesn't accept ym and sdpvr
input, use manuel difference instead
                if isempty(tol)
                    r1 = abs(real(o1.value)-real(o1 r.value))<eps;</pre>
                    r2 = abs(imag(o1.value)-imag(o1 r.value))<eps;
                else
                    r1 = abs(real(o1.value)-real(o1 r.value))<tol;</pre>
                    r2 = abs(imag(o1.value)-imag(o1 r.value))<tol;
                end
            end
                R = zeros(size(o1.value), 'logical');
                R((r1 + r2) == 2) = true;
                r = all(R, 'all');
        end
        function tol = tolReal(o1,tolstart,tolTol)
            % tolReal - Determines the highest tolerance for which the input
is real.
            % Syntax: tol = tolReal(o1, tolstart, tolTol)
            % Inputs:
                ol - The input object to be checked for reality.
                 tolstart - (Optional) The starting tolerance value for the
binary search. Default is 100.
                tolTol - (Optional) The tolerance for the binary search
convergence. Default is 1e-20.
            % Outputs:
                tol - The highest tolerance for which the input is real
through function isreal(o1, tol).
            % Example:
                tol = tolReal(o1, 100, 1e-20);
            % See also: isreal
            arguments
                01
                tolstart=100
```

```
tolTol=1e-20;
            end
            % Initialize variables
            low = 0;
            high = tolstart;
            % Binary search for minimum tolerance
            while (high - low) > tolTol
                mid = (low + high) / 2;
                if isreal(o1, mid)
                    high = mid;
                else
                     low = mid;
                end
            end
            % Return the minimum tolerance
            tol = high;
        end
        function [r,R] = isimag(o1,tol)
            \mbox{\ensuremath{\$}} ISIMAG Check if the real part of the time realization of a
PhasorArray is negligible.
               [r, R] = ISIMAG(o1, to1) checks whether the real part of the
time-domain
                realization of the PhasorArray `o1` is negligible relative
to a given tolerance.
            응
                Inputs:
            90
                    ol - (PhasorArray) The PhasorArray object to be checked.
                    tol - (double, optional) Tolerance threshold for
comparison.
                           - Default: `[]`, meaning it uses machine precision.
            응
            응
                Outputs:
                    r - (logical) True if the entire PhasorArray is purely
imaginary.
                    R - (logical array) Element-wise result indicating if
each entry is purely imaginary.
               Example:
                    [r, R] = isimag(o1, 1e-6);
            응
                See also: mimag, isreal, abs
            arguments
                ^{\circ}1
                tol=[]
            end
            o1 i = mimag(o1);
            %if o1.value is not sym, sdpvar nor ndsdpvar
            if ~(isa(o1.value, "sym") || isa(o1.value, "ndsdpvar") ||
isa(o1.value, "sdpvar"))
                if isempty(tol)
```

```
r1 = ismembertol(real(o1.value), real(o1 i.value));
                    r2 = ismembertol(imag(o1.value),imag(o1 i.value));
                else
                    r1 = ismembertol(real(o1.value), real(o1 i.value), tol);
                    r2 = ismembertol(imag(o1.value),imag(o1 i.value),tol);
                end
            else %in this case ismembertol doesn't accept ym and sdpvr
input, use manuel difference instead
                if isempty(tol)
                    r1 = abs(real(o1.value)-real(o1 i.value))<eps;</pre>
                    r2 = abs(imag(o1.value)-imag(o1 i.value))<eps;
                else
                    r1 = abs(real(o1.value)-real(o1 i.value))<tol;
                    r2 = abs(imag(o1.value)-imag(o1 i.value))<tol;
            end
                R = zeros(size(o1.value), 'logical');
                R((r1 + r2) == 2) = true;
                r = all(R, 'all');
        end
        function [r,R] = iscomplex(o1,tol)
            % ISCOMPLEX Check if the PhasorArray contains significant
imaginary components.
            응
                [r, R] = ISCOMPLEX(o1, tol) determines whether the
            응
PhasorArray `o1`
                contains significant imaginary components relative to a
given tolerance.
                Inputs:
                   o1 - (PhasorArray) The PhasorArray object to be checked.
                   tol - (double, optional) Tolerance threshold for
comparison.
                          - Default: `eps` (machine precision).
            응
               Outputs:
                    r - (logical) True if the entire PhasorArray has
significant imaginary components.
                    R - (logical array) Element-wise result indicating if
            응
each entry is complex.
            응
            응
                Example:
                    [r, R] = iscomplex(o1, 1e-6);
                See also: isreal, isimag, abs
            arguments
                01
                tol=eps
            end
            [r,R] = isreal (o1, tol);
            r = \sim r;
            R=\sim R;
        end
```

```
function [r,R] = iszero(o1,tol)
            % ISZERO Check if the PhasorArray is numerically zero.
            응
                [r, R] = ISZERO(o1, tol) determines whether the PhasorArray
`01`
                is numerically zero, based on an absolute tolerance.
            응
            용
            응
               Inputs:
                    ol - (PhasorArray) The PhasorArray object to be checked.
                    tol - (double, optional) Tolerance threshold for
comparison.
                          - Default: `0`, meaning exact zero comparison.
            응
               Outputs:
                    r - (logical) True if all entries in the PhasorArray are
within the tolerance of zero.
                    R - (logical array) Element-wise result indicating if
each entry is within tolerance.
            응
               Example:
                    [r, R] = iszero(o1, 1e-6);
               See also: pagenorm, isreal, iscomplex
            arguments
                01
                tol = 0
            end
            R = pagenorm(o1.value, "inf") <=tol;</pre>
            r = ((nnz(R)) == (2*o1.h+1));
        end
        function r = ImagRealForm(o1)
            % IMAGREALFORM Convert PhasorArray to Imaginary-Real form.
                r = IMAGREALFORM(o1) converts the PhasorArray `o1` into the
               Imaginary-Real representation, ensuring:
                    - **Negative harmonics (`-k`) store the imaginary part
(left side).**
                    - **Positive harmonics (`+k`) store the real part (right
side).**
                    - The order `k` and `-k` remain **symmetric** around the
DC component.
                This format assumes that `o1` represents a **real-valued
periodic matrix**,
            응
               meaning X - k = conj(X k).
            응
            응
                Input:
            응
                    o1 - (PhasorArray) The input PhasorArray object.
            응
            응
               Output:
                    r - (array) The Imaginary-Real form representation:
            응
                        - Dimensions: (m \times n \times (2h+1))
```

```
- Order `-k` maps to `(:,:,h-k+1)` (imaginary part,
negated).
            응
                         - Order \dot{} +k \dot{} maps to \dot{} (:,:,h+k+1) \dot{} (real part).
            응
            응
                Example:
            응
                   r = ImagRealForm(o1);
            9
                See also: RealImagForm, SinCosForm
            h=o1.h;
            o1=o1.value;
            r=cat(3,-imag(ol(:,:,l:h)),real(ol(:,:,(h+1):end)));
        end
        function r = RealImagForm(o1)
            % REALIMAGFORM Convert PhasorArray to Real-Imaginary form.
                r = REALIMAGFORM(o1) converts the PhasorArray `o1` into the
                Real-Imaginary representation, ensuring:
                    - **Negative harmonics (`-k`) store the real part (left
side).**
                    - **Positive harmonics (`+k`) store the imaginary part
(right side).**
                    - The order `k` and `-k` remain **symmetric** around the
DC component.
            응
                This format assumes that `o1` represents a **real-valued
periodic matrix**,
            응
                meaning X - k = conj(X k).
            응
            응
                Input:
            응
                    ol - (PhasorArray) The input PhasorArray object.
            용
            응
                Output:
            응
                    r - (array) The Real-Imaginary form representation:
                        - Dimensions: (m \times n \times (2h+1))
            응
                         - Order \ -k\ maps to \ (:,:,h-k+1)\  (real part).
            응
                         - Order `+k` maps to `(:,:,h+k+1)` (imaginary part).
            응
            응
            양
                Example:
            응
                    r = RealImagForm(o1);
                See also: ImagRealForm, SinCosForm
            r = flip(ImagRealForm(o1), 3);
        end
        function r = SinCosForm(o1,isReal,realTol)
            % SINCOSFORM Convert PhasorArray to Sine-Cosine form.
                r = SINCOSFORM(o1, isReal, realTo1) converts the PhasorArray
`o1`
                into the **Sine-Cosine representation**, ensuring:
                    - **Negative harmonics (`-k`) store sine coefficients
(left side).**
                    - **Positive harmonics (`+k`) store cosine coefficients
(right side).**
                    - The order `k` and `-k` remain **symmetric** around the
```

```
DC component.
            응
                Inputs:
                           - (PhasorArray) The input PhasorArray object.
                   01
            응
                    isReal - (logical, optional) If `true`, forces real-
valued output.
                              Default: `isreal(o1)`.
                   realTol - (double, optional) Tolerance for enforcing
real values.
                              Default: `1e-14`.
            응
            응
               Output:
                    r - (array) The Sine-Cosine form representation:
                        - Dimensions: (m \times n \times (2h+1))
                        - Order `-k` maps to `(:,:,h-k+1)` (sine
            응
coefficient, negated).
                        - Order `+k` maps to `(:,:,h+k+1)` (cosine
coefficient).
            응
               Example:
                  r = SinCosForm(o1, true, 1e-14);
               See also: CosSinForm, ImagRealForm
            arguments
                01
                isReal logical = isreal(o1)
                realTol = 1e-14
            end
            %normal sin / cos form
            h=o1.h;
            o1=o1.value;
            r=cat(3,1i*(flip(ol(:,:,h+2:end),3)-ol(:,:,1:h)),ol(:,:,h+1),
(o1(:,:,h+2:end)+flip(o1(:,:,1:h),3)));
            if isReal
                er = real(r-real(r));
                if norm(double(er),'fro') > realTol
                    error ("SinCosForm : Imaginary part of matrix %d is
superior to tolerance %d, switch is Real to false or adjust
tolerance", norm(double(er), 'fro'), realTol)
                end
                r=real(r);
            end
        end
        function r = AngleAmpForm(o1)
            % ANGLEAMPFORM Convert PhasorArray to Angle-Amplitude form.
                r = ANGLEAMPFORm(o1) converts the PhasorArray `o1` into
                the **Angle-Amplitude representation**, ensuring:
                    - **Negative harmonics (`-k`) store phase angles (left
side).**
            응
                    - **Positive harmonics (`+k`) store amplitudes (right
side).**
                    - The order `k` and `-k` remain **symmetric** around the
DC component.
```

```
This format assumes `o1` represents a **real-valued**
periodic matrix.
            응
                Input:
                    ol - (PhasorArray) The input PhasorArray object.
            용
            응
                Output:
            90
                    r - (array) The Angle-Amplitude form representation:
                        - Dimensions: (m \times n \times (2h+1))
            응
                        - Order \dot{}-k maps to \dot{}(:,:,h-k+1) \dot{} (phase angle).
                        - Order `+k` maps to `(:,:,h+k+1)` (amplitude).
            응
            응
            응
               Example:
                 r = AngleAmpForm(o1);
                See also: SinCosForm, ImagRealForm
            h=o1.h;
r=cat(3, angle(flip(ol.phas(1:h))), real(ol.phas(0)), abs(ol.phas(1:h)));
        end
        function r = CosSinForm(o1)
            % COSSINFORM Convert PhasorArray to Cosine-Sine form.
            % r = COSSINFORM(o1) converts the PhasorArray `o1` into
              the **Cosine-Sine representation**, ensuring:
                    - **Negative harmonics (`-k`) store cosine coefficients
(left side).**
                    - **Positive harmonics (`+k`) store sine coefficients
(right side). **
            응
                Input:
            %
                    ol - (PhasorArray) The input PhasorArray object.
            응
            응
               Output:
            응
                    r - (array) The Cosine-Sine form representation.
            응
            응
               Example:
            응
                    r = CosSinForm(o1);
                See also: SinCosForm, ImagRealForm
            r = flip(SinCosForm(o1), 3);
        end
        function r = squeeval(o1)% SQUEEVAL Evaluate and squeeze PhasorArray.
                r = SQUEEVAL(o1) evaluates the PhasorArray `o1` and applies
`squeeze`
                to remove singleton dimensions. This is useful for cases
where `o1`
            응
                represents a scalar-valued PhasorArray.
            응
            응
                Input:
                    ol - (PhasorArray) The input PhasorArray object.
```

```
응
               Output:
                r - (array) The squeezed numerical evaluation of
`o1.value`.
            응
               Example:
            응
                   r = squeeval(o1);
                See also: value, squeeze
            r=squeeze(value(o1));
        end
        function r = expandBase(o1, m)
            \mbox{\ensuremath{\$}} EXPANDBASE Insert zeroed phasors to change the frequency base.
                r = EXPANDBASE(o1, m) modifies the PhasorArray `o1` by
inserting `m-1`
               zero-valued phasors between each existing phasor,
effectively changing
               the frequency base from `w0` to `w1 = w0 / m`.
              This transformation results in a PhasorArray with an
increased harmonic
                resolution, suitable for frequency resampling or down-
sampling applications.
            응
               Input:
                 ol - (PhasorArray) The input PhasorArray.
                    m - (integer) The expansion factor (must be \geq 1).
            % Output:
                    r - (PhasorArray) The modified PhasorArray with
inserted zero phasors.
               Behavior:
                    - The harmonics are redistributed such that phasors
appear only at
                      indices `k*m`, while the in-between phasors are zeroed.
                    - The resulting PhasorArray corresponds to the same
            응
periodic function,
                     but expressed in a lower fundamental frequency `w1 =
w0/m.
            응
              Example:
                   A new = expandBase(A, 3); % Adjusts A for frequency w0/3
               See also: squishBase, pad
            h = o1.h;
            nx = size(o1,1);
            ny = size(o1, 2);
            r = zeros(nx, ny, 2*m*h+1);
            r(:,:,1:m:end) = o1.value;
            r = PhasorArray(r);
```

```
end
        function r = squishBase(o1, m)
            % SQUISHBASE Remove phasors to change the frequency base.
                r = SQUISHBASE(o1, m) modifies the PhasorArray `o1` by
            응
retaining only
                every `m`-th phasor, effectively changing the frequency base
from `w0`
               to w1 = w0 * m.
            응
                This transformation results in a reduced harmonic
resolution, increasing
                the fundamental frequency but potentially losing information
in the process.
                Input:
                    o1 - (PhasorArray) The input PhasorArray.
                    m - (integer) The reduction factor (must be \geq 1).
            90
               Output:
                    r - (PhasorArray) The modified PhasorArray with
retained harmonics.
               Behavior:
                    - Only harmonics at indices `k*m` are retained, while
            응
others are discarded.
                    - If `ol.h` is not a multiple of `m`, it is first padded
with zeros.
                    - A warning is issued if nonzero harmonics are removed,
as this can
                     lead to information loss and an inaccurate
representation.
            응
                Example:
            응
                    A new = squishBase(A, 3); % Adjusts A for frequency w0 *
3
                See also: expandBase, pad
            h = o1.h;
            nx = size(o1,1);
            ny = size(o1, 2);
            %ensure o1.h is a multiple of m, otherwise pad o1 with zeros
            if mod(h, m) \sim = 0
                o1 = pad(o1, m-mod(h, m));
                h = o1.h;
            end
            r = o1(:,:,1:m:end);
```

%evaluate the norm of deleted phasors in original phasorArray

r = PhasorArray(r);

```
deleted = o1(:,:,setdiff(1:(2*h+1),1:m:(2*h+1)));
            normDeleted = norm(deleted(:),'fro');
            if normDeleted > 1e-10
                warning ("squishBase: deleted phasors have a norm of %d, the
new phasorArray may not accurately represent the original
signal", normDeleted)
            end
        end
    end
    methods (Access=protected)
        function varargout = parenReference(obj,indexOp)
            obj.Phasor3D = obj.Phasor3D.(indexOp(1));
            if isscalar(indexOp)
                varargout{1} = pvalue(obj);
                return;
            end
            % Syntax for forwarding indexing operations
            [varargout{1:nargout}] = obj.(indexOp(2:end));
        end
        function obj = parenAssign(obj,indexOp,varargin)
            if isscalar(indexOp)
                assert(nargin==3);
                rhs = vararqin{1};
                if isa(rhs,'PhasorArray')
                    obj.Phasor3D.(indexOp) = rhs.value;
                else
                    obj.Phasor3D.(indexOp) = rhs;
                end
                return:
            [obj.(indexOp(2:end))] = varargin{:};
        end
        function n = parenListLength(obj,indexOp,ctx)
            if numel(indexOp) <= 2</pre>
                n = 1;
                return;
            end
            containedObj = obj.(indexOp(1:2));
            n = listLength(containedObj,indexOp(3:end),ctx);
        end
        function obj = parenDelete(obj,indexOp)
            obj.Phasor3D.(indexOp) = [];
        end
        function varargout=braceReference(obj,indexOp)
            if numel(indexOp(1).Indices) == 3
                argou=obj.sub(indexOp(1).Indices{1:2});
                varargout{1}=argou.phas(indexOp(1).Indices{3});
            else
```

```
if ~isscalar(indexOp)
                     for ii=1:numel(varargout)
                         varargout{ii}=varargout{ii}.(indexOp(2:end));
                     end
                end
            end
            응
                           indexOp
                           isscalar(indexOp)
            양
                           indexOp(1)
            응
                           indexOp.Type=Paren
            응
                           a=index0p
            응
                           contobj=obj.Phasor3D
                           contobj=contobj(:,:,indexOp(1))
            응
            응 응
                             obj.Phasor3D = obj.Phasor3D.(:,:,indexOp(1));
            응
                           if isscalar(indexOp)
            응
                               varargout{1} = contobj;
            용
                               return;
            응
                           end
            응
                           % Syntax for forwarding indexing operations
            응 응
                             [varargout{1:nargout}] = obj.(indexOp(2:end));
            응 응
                             objv=obj.value;
            응
                           [varargout{1:nargout}] = objv(:,:,indexOp);
        end
        function n = braceListLength(obj,indexOp,indexContext)
            if numel(indexOp) <= 2</pre>
                n = 1;
                return;
            end
            n=1;
        end
        function obj = braceAssign(obj,indexOp,varargin)
            h1=size(obj,3);
            h2=size(varargin{:},3);
            m1=size(obj,1);
            m2=size(obj,2);
            if numel(indexOp.Indices) == 3
                n1=indexOp.Indices{1};
                n2=indexOp.Indices{2};
                n3=indexOp.Indices{3};
                if ischar(n3) %on a reçu l'argument ':', donc forcement un
phasorArray
                else
                     if max(abs(n3))>obj.h
                         obj=obj.pad(max(abs(n3))-obj.h);
                     end
```

[varargout{1:nargout}]=obj.sub(indexOp(1).Indices{:});

```
n3=n3+1+obj.h;
                end
                obj(n1, n2, n3) = varargin{:};
                return
            end
            if h1<h2
                obj=obj.pad((h2-h1)/2);
            else
                varargin{:}=PhasorArray(PhasorArrayPad(varargin{:}, (h1-h2)/
2));
            end
            if isscalar(indexOp)
                switch numel(indexOp.Indices)
                     case 1
                         m=indexOp.Indices{1};
                         if numel(m) == obj.numelt %numelt renvoit la dim de
la matrice tempo ie n1 \times n2, donc ce serait un logical array en entrée
                             if any((m \sim = 0) & (m \sim = 1)) %on verifie que
c'est un logical array
                                 error('Array indices must be positive
integers or logical values.')
                             end
                             n1=find(indexOp.Indices{1});
                         else
                             n1=indexOp.Indices{1};
                         end
                         n2=1;
                         % obj
                         obj=obj{:};
                     case 2
                         n1=indexOp.Indices{1};
                         n2=indexOp.Indices{2};
                end
                 if isscalar(varargin{:})
                     varargin{:}=varargin{:}*ones(numel(n1),numel(n2));
                end
                                   obj(n1,n2,:) = vect(varargin\{:\});
                obj(n1,n2,:) = (varargin\{:\});
                obj=reshape(obj,m1,m2,[]);
                return;
            end
                           [braceReference(obj,indexOp)] = varargin{:};
        end
    end
    methods (Access=public)
        function ind = end(obj,k,n)
            s1 = size(obj, 1);
```

```
s2 = size(obj, 2);
            s3 = size(obj,3);
            sz = [s1 \ s2 \ s3];
            if n == 1
                ind = s1*s2;
                return
            end
            ind = sz(k);
        end
        function out = value(obj)
            out = obj.Phasor3D;
        end
        function out = sum(obj,dim)
            %SUM Compute the sum along a given dimension.
              out = SUM(obj, dim) returns the sum along the specified
dimension.
                If dim is 1 or 2, the output remains a PhasorArray.
            arguments
                obj
                dim=1
            end
            out = sum(obj.Phasor3D,dim);
            if dim==1 || dim==2
                out=PhasorArray(out);
            end
        end
        function out = cat(dim, varargin)
            %CAT Concatenate multiple PhasorArrays along a given dimension.
                out = CAT(dim, A1, A2, ...) concatenates multiple
            응
PhasorArrays or
               numeric arrays along the specified dimension.
            numCatArrays = nargin-1;
            if dim==1 || dim==2
                varargin2=PhasorUnif(varargin{:});
            else
                varargin2=varargin;
            end
            newArgs = cell(numCatArrays,1);
            for ix = 1:numCatArrays
                if isa(varargin2{ix},'PhasorArray')
                    newArgs{ix} = varargin2{ix}.Phasor3D;
                    newArgs{ix} = varargin2{ix};
                end
            end
            out = PhasorArray(cat(dim, newArgs{:}));
        end
```

```
function varargout = size(obj,varargin)
            [varargout{1:nargout}] = size(obj.Phasor3D, varargin{:});
        end
        function r = sdpval(o1)
            %SDPVAL Extract the numerical value of a PhasorArray whose value
is an SDPVAR.
                r = sdpval(o1) extract the numerical value of o1 a
phasorArray with NDSDPVAR 3DArray.
                Basically perform PhasorArray(value(value(o1)))
            if isa(o1.value,'ndsdpvar')||isa(o1.value,'sdpvar')||
isa(o1.value, 'sym')
                o1=o1.value;
            end
            r=PhasorArray(value(o1));
        end
        [K,P,res] = place(A,B,poles,varg) ;
        out = expm(A, varg)
        out = logm(A, varg)
    end
    methods
        function r = Value(obj)
            r=obj.value;
        end
    end
    methods (Static, Access=public)
        function out = time2Phasor(At,nT,t,varg)
            %TIME2PHASOR Convert a time-dependent matrix into a PhasorArray
representation.
            응
                out = TIME2PHASOR(At, nT, t, varg) converts a 3D array
representing a
               real, time-dependent matrix (with time stored along the 3rd
dimension)
                into a PhasorArray by computing its Fourier coefficients.
            응
            응
               Assumptions:
                 - The sampling is uniform: Ts = t(2) - t(1).
                  - The final time matches a full number of periods: t(end)
= nT*T - dt.
            응
               Inputs:
                 Αt
                           - (3D array) Time-dependent matrix, stored as
MxNxtime.
            용
                 nΤ
                           - (integer, default: 1) Number of periods
captured in `t`.
                           - (vector, optional) Time vector. If empty, it is
                  t
inferred.
```

```
varg - (Optional) Name-value pair arguments:
                     'truncIndex' (integer, default: Inf) - Maximum
harmonic order for truncation.
                                  (logical, default: true) - Enforce
           응
                     'real'
conjugate symmetry (A k = conj(A - k)).
                                (integer, default: 3) - Dimension
                     'timeDim'
along which time varies.
           응
              Outputs:
                 out - (PhasorArray) The computed PhasorArray
            응
representation of `At`.
            응
            응
               Example:
            응
                 t = linspace(0, 2*pi, 100);
                 At = cat(3, sin(t), cos(t));
            응
                Ph = time2Phasor(At, 1, t, 'truncIndex', 5);
              See also: PhasorArray, TimeArray2Phasors
           arguments
           Αt
           nT=1
           t=[]
           varg.truncIndex=Inf
           varg.real=true
           varq.timeDim=3
           end
PhasorArray(TimeArray2Phasors(At,nT,t,'truncIndex',varg.truncIndex,'isReal',v
arg.real, 'timeDim', varg.timeDim));
       end
        function [phasorArr,feval] = funcToPhasorArray(func, T, n,varg)
            %FUNCTOPHASORARRAY Convert a time function into a PhasorArray.
               [phasorArr, feval] = FUNCTOPHASORARRAY(func, T, n, varg)
evaluates a
              time-dependent function over a uniform time grid and
converts it into
              a PhasorArray using a Fourier transform.
            응
               Inputs:
                 func - (function handle) Time-dependent function, f(t),
returning an MxN matrix.
                 T - (scalar, default: 1) Period of the function.
                      - (integer, default: 4) Number of frequency bins used
                 n
(2^n time steps).
                varg - (Optional) Name-value pair arguments:
                     'reduce'
                                 (logical, default: true) - Reduce the
output PhasorArray.
                     'reduceTol' (double, default: 1e-15) - Threshold for
           응
reducing small harmonics.
              Outputs:
```

```
phasorArr - (PhasorArray) The computed PhasorArray
representation of `func(t)`.
                  feval
                           - (struct) Structure containing function
            응
evaluations for debugging:
                      - feval.T : Period used.
            응
                      - feval.n : Number of frequency bins.
            용
                      - feval.dt : Time step.
                      - feval.func : Function handle used.
                      - feval.At : Evaluated matrix over time.
                      - feval.t : Time vector used.
            응
            응
                Example:
                 func = Q(t) [sin(t); cos(t)];
            응
            응
                  T = 2*pi;
            응
                 n = 5;
            응
                 [Ph, fData] = funcToPhasorArray(func, T, n);
                See also: time2Phasor, PhasorArray
            arguments
                func (1,1) {mustBeA(func, 'function handle')} = @(t)
ones (3,3)
                T (1,1) {mustBeNumeric, mustBePositive} = 1
                n (1,1) {mustBeNumeric, mustBePositive, mustBeInteger} = 4
                varq.reduce = true;
                varg.reduceTol = 1e-15;
            end
            % Calculate dt
            dt = T / (2^n);
            % Create a time vector
            t = 0:dt:T-dt;
            % Evaluate the function over the time vector
            At = arrayfun(func, t, 'UniformOutput', false);
            At = cat(3, At\{:\});
            AtT = func(T);
            %evaluate the jump between the last and first value
            if norm(At(:,:,end)-At(:,:,1),'fro')>1e-10
                warning ('The function has discontinuities, or a steep
derivative at the end, or has a jump between the last and first value, the
result may be incorrect. Jump value is %d (froebenius norm of f(T-dt)-
f(0))', norm(At(:,:,end)-At(:,:,1),'fro'));
            end
            if norm(At(:,:,1)-AtT,'fro')>1e-10
                warning('A(T) is different from A(0), the result may be
incorrect. Jump value is %d (froebenius norm of f(T)-f(0)), resulting
phasorArray is a periodic function with jump at time T', norm(At(:,:,1)-
AtT, 'fro'));
            end
```

```
% Convert the time-domain function values to a PhasorArray
            phasorArr = PhasorArray.time2Phasor(At);
            if varq.reduce
                phasorArr = reduce(phasorArr,
"reduceThreshold", varg.reduceTol, "reduceMethod", "relative", "excludeOPhasor", f
alse, "hardThresholdPhasors", true);
            end
            if nargout>1
                feval=struct;
                feval.T=T;
                feval.n=n;
                feval.dt=dt;
                feval.func=func;
                feval.At=At;
                feval.t=t;
                feval.plot = @(n,alone) At2plot(n,alone);
            end
            function At2plot(n,alone)
                %PLOTFUNC Plot time-domain representation of function
evaluations.
                    plotFunc(numPeriods, mode) plots the function
evaluations over `numPeriods`
                % periods. The `mode` argument can be 'alone' (plot time-
domain only) or 'both'
                    (overlay with the PhasorArray reconstruction).
                응
                응
                응
                   Example:
                     feval.plot(4, 'both');
                응
                if nargin==0
                    n=4;
                    alone="alone";
                end
                if nargin==1
                    if (ischar(n) || isstring(n))
                        alone=n;
                        n=4;
                    elseif nargin==1 && isnumeric(n)
                        alone="alone";
                    end
                end
                %assert if alone is "alone" or "both"
                assert(strcmp(alone, "alone") ||strcmp(alone, "both"), "alone
must be 'alone' or 'both'");
```

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```
dt=T/(100);
                t = 0:dt:n*T;
                At = arrayfun(func, t, 'UniformOutput', false);
                At = cat(3, At\{:\});
                %plot each A(i,j,t) as a function of time
                for i=1:size(At,1)
                    for j=1:size(At,2)
                        subplot (size (At, 1), size (At, 2), (i-1)*size (At, 2)+j);
                        plot(t, squeeze(At(i,j,:)));
                        title(sprintf('A {%d,%d}(t)',i,j));
                    end
                end
                if strcmp(alone, "both")
                    hold on;
                    phasorArr.plot(T,[0 n*T]);
                    hold off;
                end
            end
        end
        function out = cqt2ScalarPhasor(cqtobj,varg)
            %CQT2SCALARPHASOR Convert a CQT object to a ScalarPhasorArray
representation.
            응
            응
                out = CQT2SCALARPHASOR(cgtobj, varg) extracts the symbolic
Toeplitz
                representation of a compact quasi-Toeplitz (CQT) object and
            응
maps it
            용
               to a ScalarPhasorArray.
            응
                Inputs:
                  cqtobj - (CQT object) The input compact quasi-Toeplitz
object.
            응
                 varg - (Optional) Name-value pair arguments:
                      'isReal' (logical, default: false) - If true, enforces
real-valued output
                                                            by ensuring
conjugate symmetry in
                                                             the phasor
representation.
                Outputs:
                  out - (ScalarPhasorArray) The extracted and converted
phasor representation.
            응
            응
                Example:
            응
                  cqtA = cqt(rand(10), rand(10)); % Example CQT matrix
            응
                  phasorA = cqt2ScalarPhasor(cqtA, 'isReal', true);
            응
                See also: ScalarPhasorArray, cqt
            arguments
```

```
cqtobj %object of class cqt
                varg.isReal = false %enforce real output (ie conjugate
phasore pos/neq)
            end
            [nsTA, psTA] = symbol (cqtobj);
            n=min(numel(nsTA), numel(psTA));
            nsTA=nsTA(1:n);
            psTA=psTA(1:n);
            out=ScalarPhasorArray([flip(nsTA(2:end)) psTA]);
            if varq.isReal
                out = mreal(out);
            end
        end
        function obj = empty(varargin)
            %EMPTY Create an empty PhasorArray of specified size.
               See also: zeros, ones, eye
            obj = PhasorArray(double.empty(varargin{:}));
        end
        function obj = zeros(vararqin)
            %ZEROS Create a zero-filled PhasorArray of specified size.
            % See also: ones, eye, empty
            obj = PhasorArray(zeros(varargin{:}));
        end
        function obj = ones(varargin)
            %ONES Create a PhasorArray filled with ones.
               See also: zeros, eye, empty
            obj = PhasorArray(ones(varargin{:}));
        end
        function obj = eye(varargin)
            %EYE Create a PhasorArray with identity matrices at all
harmonics.
              EYE(n) returns an n \times n \times 1 PhasorArray filled with
identity matrices.
                EYE (n, m) returns an n \times m \times 1 PhasorArray filled with
            용
identity matrices.
            % EYE(n, m, h) returns an `n × m × (2h+1)` PhasorArray where
each slice along the 3rd dimension is an identity matrix.
            % Unlike a standard identity matrix, this function replicates
the identity structure across all harmonic indices.
            % See also: zeros, ones, empty
            if nargin==1
                varargin{2}=varargin{1};
                varargin{3}=1;
            elseif nargin==2
                varargin{3}=1;
            elseif nargin==3
            else
            u=repmat(eye(varargin{1:2}),[1 1 2*varargin{3}+1]);
            obj = PhasorArray(u);
        end
```

```
function obj = randomPhasorArrayWithPole(nx,poles,T,varg)
            % RANDOMPHASORARRAYWITHPOLE Generate a PhasorArray with
prescribed poles.
            응
            응
                OBJ = RANDOMPHASORARRAYWITHPOLE(NX, POLES, T, VARG)
constructs a
                time-periodic matrix `A(t)` with the specified eigenvalues
(`poles`).
                A first random matrix \hat{A} is generated, and a B*G term is
computed to
                ensure the desired eigenvalues by performing pole placement.
            응
                The resulting PhasorArray `OBJ` is the difference `A-BK`,
where `K` is
               the obtained feedback gain.
            응
            응
                Inputs:
            응
                 NX
                        - (integer) The size of the square matrix.
                  POLES - (vector) Desired eigenvalues for `A(t)`, must have
length NX.
                        - (scalar, optional) Period of the matrix. Default
                  Т
is `2*pi`.
            응
                  VARG - (optional) Name-value pair arguments:
                            - 'h'
                                        (integer) Harmonic order (default:
`5`).
                            - 'BG'
                                        (PhasorArray) Custom B*G term
(default: random).
                            - 'isReal' (logical) Force a real-valued matrix
(default: `true`).
            응
            응
                Outputs:
                  OBJ - (PhasorArray) Generated PhasorArray with specified
poles.
                See also: Sylv harmonique, random
            arguments
                nx
                poles
                T = 2*pi
                varg.h = []
                varg.BG = []
                varg.isReal = true
            end
            if isempty(varg.h) && isempty(varg.BG)
                varq.h=5;
                A = PhasorArray.random(nx,nx,varg.h);
                BG = PhasorArray.random(nx,nx,varg.h);
            elseif isempty(varg.BG)
                A = PhasorArray.random(nx,nx,varg.h);
                BG = PhasorArray.random(nx,nx,varg.h);
            elseif isempty(varg.h)
                varg.h = varg.BG.h;
                A = PhasorArray.random(nx,nx,varq.h);
            else
                A = PhasorArray.random(nx,nx,varg.h);
```

```
end
            assert(numel(poles) ==nx)
            La = PhasorArray(diag(poles));
            %Solve the appropriate Sylvester equation
            P = PhasorArray(Sylv harmonique(-A, La, BG, 4*varg.h, 2*pi/T));
            %Compute K
            K = BG/P;
            %compute the new A with appropriate eigen values
            obj = A-K;
        end
        function obj = random(nx,ny,h,arg)
            % RANDOM Generate a random PhasorArray with optional structure
constraints.
                OBJ = RANDOM(NX, NY, H, ARG) generates a random 3D
PhasorArray representing
                a time-periodic matrix with harmonics and structural
constraints.
            응
                Inputs:
                 NX - (integer) Number of rows.
                  NY - (integer) Number of columns.
                 H - (integer) Number of harmonics.
            응
                  ARG - (optional) Name-value pair arguments:
                          - 'time structure' (char) Time-domain structure:
                                 'real', 'symmetric', 'antisymmetric', 'sdp',
'hermitian',
                                 'cmplx', 'retroHermitian', etc. (default:
'real').
                          - 'hurwitzeig' (vector) Predefined Hurwitz
eigenvalues (default: `[]`).
                          - 'T'
                                           (scalar) Time period (default:
`1`).
                          - '0'
                                           (matrix) Optional quadratic matrix
(default: `[]`).
                          - 'average power decay' (scalar) Power decay rate
of harmonics (default: `2`).
            응
            응
               Outputs:
            응
                 OBJ - (PhasorArray) Randomly generated PhasorArray.
                See also: rand phasor
            arguments
                nx
                ny
                arg.time structure {mustBeMember(arg.time structure,
{'real', 'symetric', 'antisymetric', 'sdp', 'hurwitz', 'Q-
spec','hermitian','cmplx','retroHermitian'})} = 'real'
                arg.hurwitzeig=[]
                arg.T=1
                arg.Q=[]
```

```
arg.average power decay=2;
            end
            if nargin ==1
                switch numel(nx)
                    case 1
                    case 2
                        ny = nx(2);
                        nx = nx(1);
                        h = 0;
                    case 3
                        ny = nx(2);
                        h = nx(3);
                        nx = nx(1);
                end
            end
            arg.output ='PhasorArray'; %or PhasorArray
            if ~isempty(arg.hurwitzeig) && strcmp(arg.time structure, 'real')
                arg.time structure = 'hurwitz';
            end
            C=namedargs2cell(arg);
            obj = rand phasor(nx,ny,h,C{:});
        function A = sym(nx,ny,h,name)
            % SYM Construct a symbolic PhasorArray with structured harmonic
components.
            응
                A = SYM(NX, NY, H, NAME) creates an NX-by-NY PhasorArray
where each
                element is a symbolic expression representing harmonic
components.
            응
            응
                Inputs:
            응
                       - (integer) Number of rows (default: 1).
                  NX
                  NY - (integer) Number of columns (default: NX).
                       - (integer) Number of harmonics (default: 0).
                  NAME - (string or cell) Base name for symbolic variables
(default: "a").
            응
            응
                Outputs:
            응
                 A - (PhasorArray) Symbolic PhasorArray with structured
harmonics.
            응
            응
                Notes:
                  - Each entry in A has a symbolic expression for phasors:
                    `A 0`, `A plus k`, `A minus k` for k > 0.
                  - If NAME is a cell array, it must match the size of NX \times
NY.
                  - For scalar NX \times NY, NAME is used as the base for all
symbols.
                See also: sym, ScalarPhasorArray, PhasorArray.
            arguments
```

```
nx=1
                ny=nx
                h = 0
                name ="a"
            end
            if (nargin == 1) || (nargin==2 && (ischar(ny) || isstring(ny) ||
iscell(ny)))
                    if ischar(ny) || isstring(ny) || iscell(ny)
                        name = ny;
                    end
                    switch numel(nx)
                        case 1
                            ny = nx;
                            h = 0;
                        case 2
                            ny = nx(2);
                            nx = nx(1);
                            h = 0;
                        case 3
                            ny = nx(2);
                            h = nx(3);
                            nx = nx(1);
                        otherwise
                            error('cannot specify a dimension input of
length >3')
                    end
            end
            if max(ny,nx) == 1
                name={name};
            end
            if iscell(name)
                assert(numel(name) ==nx*ny);
                A=PhasorArray.zeros(nx,ny);
                A.Phasor3D = sym(A.Phasor3D);
                for ii = 1:numel(name)
                    clear ap am a0 a
                    name i=name{ii};
                    ap = sym(name_i+"_plus_",[1 h]);
                    am = sym(name i+" minus ",[1 h]);
                    a0 = sym(name i+" 0");
                    a = cat(2, flip(am), a0, ap);
                    A{ii} = ScalarPhasorArray(a);
                end
            else
                ap = sym(name+"__%d__%d_plus_%d",[nx ny h]);
                am = sym(name+" %d %d minus %d",[nx ny h]);
                a0 = sym(name+" %d %d 0",[nx ny]);
                a = cat(3, flip(am, 3), a0, ap);
                A= PhasorArray(a);
            end
```

```
end
        function P = ndsdpvar(n1, n2, h, varg)
            % NDSPDPVAR Construct an `sdpvar`-based PhasorArray of specified
size and structure.
               P = NDSPDPVAR(N1, N2, H, <name-value arguments>) generates
            응
an SDP variable-based
               PhasorArray suitable for optimization in YALMIP.
            응
            응
               Inputs:
           용
                N1 - (integer) First dimension size.
                N2 - (integer, optional) Second dimension size (default:
N1).
            응
                H - (integer, optional) Number of harmonics (default: 0).
            응
            % Name-Value Arguments:
                  'PhasorType' (char) - Defines the structure of the phasor
(default: 'symmetric').
                                       Options: 'symmetric', 'full',
'diagonal', etc.
                                       See YALMIP documentation for the
complete list.
                'real' (logical) - If true, ensures conjugate symmetry
for real-valued signals (default: true).
            응
            응
               Outputs:
                P - (PhasorArray) PhasorArray containing `sdpvar` elements.
            응
            응
              Notes:
                - The `PhasorType` argument defines **phasor structure**,
not time-domain structure.
            % - Example: A **Hermitian** phasor structure enforces:
                       conj(A(ij)(t)) = A(ji)(-t)
                  However, for A(t) to be Hermitian in the **time
domain**, phasors must satisfy:
                     A k = ctrans(A - k)
                 - If `real=true`, ensures the phasor array represents a
real-valued periodic matrix.
            % - If `h>0`, higher-order harmonics are created as
additional `sdpvar` variables.
              Example:
                P = PhasorArray.ndsdpvar(4,4,5, 'PhasorType', 'symmetric',
'real', true);
                   -> Produces a real-valued P(t), with 5 harmonics (size
11 along the third dimension),
                       and enforces symmetry (i.e., P ij(t) = P ji(t)).
            응
            응
                A = PhasorArray.ndsdpvar(4,4,5, 'PhasorType', 'full',
'real', true);
                   -> Produces a real-valued A(t) with no additional
structure constraints.
               See also: sdpvar, PhasorArray, PosPart2PhasorArray.
```

```
arguments
                n1
                n2=n1
                h=0
                varg.PhasorType='symmetric'
                varg.real=true
            end
            if nargin ==1
                 switch numel(n1)
                     case 1
                     case 2
                         n2 = n1(2);
                         n1 = n1(1);
                         h = 0;
                     case 3
                         n2 = n1(2);
                         h = n1(3);
                         n1 = n1(1);
                 end
            end
            if n1 \sim = n2
                 if ismember(varg.PhasorType, {'symmetric'})
                     warning('non square matrix, PhasorType switched to
"full"')
                     varg.PhasorType='full';
                 end
            end
            if varg.real
                P1=(ndsdpvar(n1, n2, 1, varg.PhasorType, 'real'));
                     P2=(ndsdpvar(n1, n2, h, varg.PhasorType, 'complex'));
                     P = PosPart2PhasorArray(P1, P2);
                 else
                     P=PhasorArray(P1);
                 end
            else
P=PhasorArray(ndsdpvar(n1,n2,2*h+1,varg.PhasorType,'complex'));
            end
        end
    end
end
ans =
  PhasorArray with properties:
    Phasor3D: 1
```

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