

Step Map and Plot (SMP) in Open Calphad (OC)

Basic description and documentation

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NOTE: This documentation unfinished and very rudimentary.

This is part of the Open Calphad (OC) documentation of the free software. The other parts are the General Thermodynamic package (GTP) the minimizer (HMS) and the software interface (OCTQ).

The software is available on <http://www.opencalphad.org> or on the opencalphad repository at <http://www.gitbub.com>.

An introduction to the Open Calphad initiative can be found in [15Sun1].

Contents

1	Introduction	4
2	Property diagrams	4
3	Phase diagrams	5
4	Basic data structures for STEP and MAP	7
4.1	Dependencies	7
4.2	The line record	7
4.3	The node record	8
4.4	The node record	9
4.5	The save record	10
4.6	The graphics option record	10
5	Subroutines and functions	11
5.1	The setup routine for step and mapping	11
5.2	Converts a start equilibrium to a start point along a line	11
5.3	Changes the variable used for mapping	12
5.4	Finding a line to follow	12
5.5	Storing a calculated point on the line	13
5.6	Storing the last equilibrium along a line	13
5.7	Changing the axis for next equilibrium calculation	13
5.8	Calculating the next equilibrium along a line	14
5.9	Calculating a node point	14
5.10	Creates a node with several lines	15
5.11	Reserves a record to save a calculated equilibrium	15
5.12	Finds a line to be calculated	16
5.13	Creates records for saveing equilibria	16
5.14	Checks if the diagram has tie-lines in the plane	16
5.15	Checks if the node point is an invariant equilibrium	16
5.16	Tries to handel problems when mapping	17

5.17 Using smaller steps when there are convergence problems	17
5.18 Lists stored equilibria	17
5.19 Plots calculated diagrams	17
5.20 Calculates equilibria for one phase at a time	18
6 Summary	18

1 Introduction

The calculation of phase diagram is an important part of any thermodynamic software. A phase diagram gives a general “map” of a system for varying external conditions. Behind the phase diagram lies many individual equilibrium calculations which determine the solubility of the different elements in the phases. This is the basis of the so called “Calphad” technique. Binary phase diagrams can be found in drawn collections and there are almost 1000 binary systems that has been assessed thermodynamically, although many of them are not using compatible models and some are not of very high quality.

For systems with 3 or more components the phase diagrams cannot be plotted as easily as a binary system although the calculations is not much more complicated using modern software. In many cases so called “property diagrams” calculated with a single variable condition, for example T , are more useful in order to understand a system at a particular composition.

At equilibrium a multicomponent thermodynamic system may consist of several phases. The Gibbs energy of each phase is modeled independently in the thermodynamic software package as a function of T, P and the phase composition. The equilibrium of a system is found by minimizing the Gibbs energy for the given set of conditions using the thermodynamic models. At equilibrium the system may consist of one or more phases with different amounts and compositions. By varying one or more conditions along an axis it is possible to generate any kind of property and phase diagrams.

To calculate a diagram the user must first provide the conditions needed to calculate a single equilibrium, a start point, and then select one, two or more of the conditions in his system as axis variables before trying to generate a diagram. The simplest set of conditions is on T, P and the overall composition. For such conditions there is always a well defined equilibrium. But the user may have conditions on the composition of a phase or the chemical potentials, μ of a component or the enthalpy of the system. In such cases the specified equilibrium may not exist.

2 Property diagrams

For a property diagram the user must select one of his conditions as axis variable and give a minimum and maximum value for the axis and an increment. The STEP procedure will start at the start point and change the axis variable with the prescribed increment until it has calculated all equilibria between the minimum and maximum axis values. The calculated equilibria will be stored in a separate data array.

Whenever there is a change in the set of stable phases the STEP procedure in the OC software will calculate the exact value of the axis variable at the phase set change. Such

a node point will be stored in the normal equilibrium list with a name like _STEP_ij where ij is a sequential index.

3 Phase diagrams

A phase diagram has at least two axis and the lines (or surfaces in 3D and hyper-surfaces in higher dimensions) in a phase diagram separate regions with different sets of stable phase. Phase diagrams in 3D may be useful for teaching ternary phase diagrams but of limited interest in general as most real alloys have more then 3 components and phase diagrams for such systems cannot be drawn in full. This section will be limited to phase diagrams calculated and plotted in 2D, the most common presentation. Along the lines in such a phase diagram we thus have zero amount of one or more phases and the lines were called “zero phase fraction”, or ZPF, lines for short by Moral[84Mor].

To obtain a 2D presentation of a multicomponent phase diagram we must make various kinds of sections or projections. For a 2D presentation of a projection some lines which seem to intersect may not be in the same plane. Thus it can be complicated to draw and understand a phase diagram but with an equilibrium software like that used in OC[15Sun2] it is easy to calculate an equilibrium with the amount of a specific phase equal to zero and then follow the line for such an equilibrium along an axis specified for the phase diagram. In fact one can follow such a line using more than 2 axis as the method to replace each axis condition, except one, with a fix phase, can be repeated.

While following a ZPF line a new phase may become stable, or an existing phase disappear, and the phase diagram software will then generate a node point and new lines with a different phase with zero amount will exit from this node point. A line may also terminate at the limit of an axis variable.

When following a line the SMP package does not use the gridminimizer as the results from the previous calculation provides good start values for the next calculation when there is just a small change in one condition. But when calculating a node point the grid minimizer should be used to test that one has not walked into a miscibility gap. Also along a line one should test regularly at each 10'th or 20'th step that there is no problem by calling the grid minimizer. But in the current version this has not yet been implemented.

As explained in any thermodynamic textbook the Gibbs phase rule determines the number of conditions needed to determine an equilibrium

$$f = n + 2 - p \tag{1}$$

where f is the degrees of freedoms, n the number of components and p the number of stable phases. To calculate an equilibrium $f = 0$ is necessary and for a unary (single component) system that means 3 conditions, for a binary 4 etc.

The type of phase diagram that is calculated depends on the number of potentials and (normalized) extensive variables used as conditions. One should also realize that the phase diagram is independent of the size of the system. We may also draw the phase diagram using other properties than those used as conditions in order to calculate the lines in the diagram.

A special type of phase diagram has “tie-lines in the plane” and in such a case a single equilibrium calculation can provide points on both lines separating this phase region. A tie-line is a line in a two-phase region connecting the compositions of the two phases in equilibrium. The most common type of phase diagram is a binary T -composition diagram where the single phase regions are separated by two-phase regions and the limiting compositions of this two-phase equilibrium are in the same plane in the diagram.

For unary diagrams the only variables are T and P , with the corresponding extensive variables S and V . When plotted with T and P as variables there are only single phase regions in the phase diagram. The lines represent the two-phase regions where one phase disappears and another become stable at the same value of T and P . However if it is plotted with V as one axis variable there are two-phase regions where the amount of the two phases varies with the axis variables.

Also in binary phase diagrams the lines are “zero phase fraction” lines but as a binary system has tie-lines in the plane most lines also represent solubility limits of a single phase region. Thus one can apply many many relations between the lines, like the lever rule, which is not generally applicable in multicomponent systems.

For multicomponent systems we normally do not have tie-lines in the plane, unless all conditions except one are potentials. But the rule is the same that the lines separate areas with different sets of stable phases. In multicomponent systems with two or more extensive variables as conditions, there is a simple rule that the number of stable phases always changes by unity for each line one cross. It can be +1 or -1 but never 2 or more. If there is an axis that is a potential, like T , and only a few elements, there is a small chance that one has an invariant equilibrium in the plane of the diagram and crossing a line representing an invariant equilibrium. Such an invariant must be at a constant value of the potential, and there is no change in the number of stable phase crossing it (except at points where other lines end at this line), but the phases stable on either side are not the same. This is the same as crossing an invariant line in a binary T-x diagram.

The reasoning here is to explain that it is quite easy to calculate a multicomponent phase diagram, we just follow all lines that represent zero amount of a phase for the given set of conditions. But it can be quite difficult to plot the diagram as will be discussed in this documentation in the next release of this software.

The mapping of a phase diagram is thus very similar to the stepping used for a property diagram. The second axis (or third axis etc) in the phase diagram has been

replaced by a fixed phase and then we just step in the other axis keeping these phases fixed. During the mapping we may have to change which axis variable that is used for stepping along the line. And of course take care to generate node points when a new phase wants to be stable or one wants to disappear because there are several ZPF lines meeting at such a point.

4 Basic data structures for STEP and MAP

4.1 Dependencies

As the step and map procedures depend on the equilibrium calculation software many of these data structures, in particular the `gtp_equilibrium_data` record type. This is described in the GTP documentation, others may be found in the HMS documentation.

4.2 The line record

This record keeps a linked list of equilibrium records calculated along a line.

```

TYPE map_line
! This is record contains a list of calculated equilibria along a line
! These are pointers to map_node records at start and end of line.
    type(map_node), pointer :: start,end
! For threading this record must be separate from other threads
! This record is created by calling calceq7 the first equilibrium of line
    type(meq_setup) :: meqrec
! the active ceq record must be stored together with all others in order to be
! copied from the ceq record saved with the node record when line starts
    type(gtp_equilibrium_data), pointer :: lineceq
! this is the number of calculated equilibria for this line and the index
! of the first and last one stored.
! The stored equilibria has an internal next link.
! lineid is a sequential index of the lines. done is negative if done
! nfixphases are the number of fixed phases replacing axis conditions
    integer number_of_equilibria,first,last,lineid,done,nfixphases
! This is used during mapping to identify lines that have the same fixed phases
! if we have 3 or more axis there can be 2 or more fix phases along the line??
    type(gtp_phasetuple), dimension(:), allocatable :: linefixph
! Save the phase tuple representing the phase fix at start node here
! If it wants to be stable at first step along a line change axis direction
!     type(gtp_phasetuple) :: nodfixph
! This is the phase index in the phr array

```

```

        integer nodfixph
! We must also save the number and set of stable phases and theit amounts
! as we will have different stable phases for different lines
        integer nstabph
        type(gtp_phasetuple), dimension(:), allocatable :: stableph
        double precision,      dimension(:), allocatable :: stablepham
! axandir is set when linenode is created to the axis and direction for first
! step from the node. It can be changed to another axis and direction
! during map and indicate the current taxis with active condition
! axchange remember the equilibrium number for an axis change
        integer axandir,axchange
! more is 1 while following the line, 0 for last equilibrium, -1 when finished
! termerr is zero unless line terminated with error
! problem is nonzero if map_problems has been called
        integer more,termerr,problems
! firstinc is a value to add the the axis variable for the first equilibrium
! to avoid finding the node point again. Evenvalue is the next value
! to calculate during a step calculation. Both set when creating the node.
! At start nodes they are zero
        double precision firstinc,evenvalue
! During map the last axis values for ALL axis are stored here
        double precision, dimension(:), allocatable :: axvals
! If tie-lines in the plane we must also check the axis valies for
! the other line as we may have to change the fix phase
        double precision, dimension(:), allocatable :: axvals2
! save previous values of axvals to handle axis changes ...
        double precision, dimension(:), allocatable :: axvalx
! factor to control length of step in axis with axtive condition
        double precision :: axfact
end TYPE map_line

```

4.3 The node record

This record contain the equilibrium at a node points where several lines meet.

```

TYPE map_node
! this record organizes the step/map results. Whenever there is a
! change of the set of stable phases a node record is created and it
! can have links to several map_line records. The map node record has a
! link to a gtp_equilibrium_data record (ceq) for the equilibrium at the node.
! This is copied to the map_line record when this is activated.
! In the map_line record an axis and direction to start is stored.

```



```

! NOTE all gtp_equilibrium_data (ceq) records are pointers to the global
! array as new composition sets may be created along any line.
! The node record is identified by the set of stable phases and the
! chemical potentials of the components. One must be able to identify the
! node as one may find the same node following different lines.
! locally stored linerecords for lines exiting the node
      type(map_line), dimension(:), allocatable :: linehead
! links to other nodes
! plotlink is used to overlay two or more map or step commands
      type(map_node), pointer :: first,next,previous,plotlink
! saved copy of the meqrec record used to calculate the node
      type(meq_setup) :: meqrec
! link to saved copy of the equilibrium record
      type(gtp_equilibrium_data), pointer :: nodeceq
! link to array of saved equilibrium record. (only maptop?)
      type(map_ceqresults), pointer :: saveceq
! copy of nodeceq in saveceq (composition sets not updated but needed for plot)
      integer savednodeceq
! type_of_node not used?? should identify invariants when tie-line not in plane
! lines are number of line records
! noofstph is number of stable phases (copied from meqrec)
! tieline_inplane is 1 if so, 0 if step, -1 if no tie-lines (only maptop)
! number_ofaxis is the number of axis, 1=step; (only maptop)
      integer type_of_node,lines,noofstph,tieline_inplane,number_ofaxis
! seqx is unique identifier for a map node
! seqy unique identifier for maplines, incremented for each line (only maptop)
      integer seqx,seqy
! nodefix is the phase held fix when calculating node (negative if removed)
      type(gtp_phasetuple) :: nodefix
! Value of T and P, copied from meqrec
      double precision, dimension(2) :: tpval
! chemical potentials, copied from meqrec
      double precision, dimension(:), allocatable :: chempots
! stable phase+compset, copied from meqrec (not used?)
      type(gtp_phasetuple), dimension(:), allocatable :: stable_phases
end TYPE map_node

```

4.4 The axis record

This record contains a description of the axis set for the mapping of the phase diagram.

```

TYPE map_axis

```

```

! description of the axis variables used for step/map
! The axis condition in bits and pieces
    integer nterm,istv,iref,iunit
    integer, dimension(:,:), allocatable :: indices
    type(gtp_state_variable), dimension(:), allocatable :: axcond
    double precision, dimension(:), allocatable :: coeffs
! the min, max and increment along the axis
    double precision axmin,axmax,axinc
! more must be initiated to 0, if nonzero replaced by a fixed phase
! seqz is the sequential index of the condition in the list (this is not
! changed if if conditions are added (at the end) or deleted (active=1)
! we cannot use a pointer as that depend on the current equilibrium.
    integer more,seqz
! This is the last succesfully calculated axis value
    double precision lastaxval
end TYPE map_axis

```

4.5 The save record

This record contains calculated equilibria.

```

TYPE map_ceqresults
! stores calculated equilibrium records
    integer size,free
    TYPE(gtp_equilibrium_data), dimension(:), allocatable :: savedceq
end TYPE map_ceqresults

```

4.6 The graphics option record

The axis and other options for plotting the diagram are stored here.

```

TYPE graphics_options
! setting options for the plotting, this replaces most arguments in the call
! to ocplot2(ndx,pltax,filename,maptop,axarr,form)
! ndx is number of plot axis, pltax is text with plotaxis variables
! filename is intermediary file (maybe not needed)
! maptop is map_node record with all results
! form is type of output (screen or postscript or jpeg)
    integer rangedefaults(3)
    double precision, dimension(3) :: plotmin,plotmax
    double precision, dimension(3) :: dfltmin,dfltmax

```

```

! labeldefaults 0 if default, 1 if text provided in plotlabels
! linetype 0 is black full line, >100 is symbols
      integer labeldefaults(3),linetype
! label 1 is heading, 2 is x-axis text, 3 is y-axis text
      character*64, dimension(3) :: plotlabels
      logical gibbstriangle
! the set key command in GNUPLOT specifies where the line id is written
! it can be on/off, placed inside/outside, left/right/center, top/bottom/center,
! and some more options that may be implemented later ...
      character labelkey*24
! many more options can easily be added when desired, linetypes etc
end TYPE graphics_options

```

5 Subroutines and functions

5.1 The setup routine for step and mapping

This routine organizes the STEP or MAP procedure.

```

      subroutine map_setup(maptop,nax,axarr,starteq)
! main map/step routine
! maptop is the main map_node record which will return all calculated lines.
! nax is the number of axis (can be just one for STEP)
! axarr is an array of records specifying the axis for the step/map
! starteq is an equilibrium data record, if there are more start equilibria
! they are linked using the ceq%next index
      implicit none
      integer nax,nsp
      type(map_axis), dimension(nax) :: axarr
      TYPE(gtp_equilibrium_data), pointer :: starteq
      TYPE(map_node), pointer :: maptop

```

5.2 Converts a start equilibrium to a start point along a line

```

      subroutine map_startpoint(maptop,nax,axarr,inactive,ceq)
! convert a start equilibrium to a start point replacing all but one axis
! conditions with fix phases. The start equilibrium must be already
! calculated. ceq is a datastructure with all relevant data for the equilibrium
! A copy of ceq and the corresponing meqrec must be made and linked from maprec
! the axis conditions replaced by fix phases are inactive
!

```

```

implicit none
TYPE(gtp_equilibrium_data), pointer :: ceq
TYPE(map_node), pointer :: maptop
integer nax
integer inactive(*)
type(map_axis), dimension(nax) :: axarr

```

5.3 Changes the variable used for mapping

If the slope of the axis variables changes this routine can change the variable used for the equilibrium calculation.

```

subroutine map_replaceaxis(meqrec,axactive,ieq,nax,axarr,tmpline,&
    inactive,ceq)
! replace an axis condition with a fix phase
    implicit none
! axactive is the axis with active condition, ieq is number of exiting lines
! ieq is the number of lines exiting from the startpoint
! nax is number of axis, axarr are description of the axis
! tmpline is to transfer some line data to calling routine
! inactive is not really used.
    type(meq_setup), pointer :: meqrec
    integer nax,axactive,ieq
    type(map_axis), dimension(nax) :: axarr
    type(gtp_equilibrium_data), pointer :: ceq
    type(map_line), dimension(2) :: tmpline
    integer inactive(*)

```

5.4 Finding a ZPF line to follow

```

subroutine map_startline(meqrec,axactive,ieq,nax,axarr,tmpline,ceq)
! find a phase to fix to replace an axis condition when we
! do not have tie-lines in the plane or when we
! have tie-lines in the plane but start in a single phase region
! meqrec is equilibrium record already initiated
! axactive is set to the axis with active condition
! ieq is the number of lines exiting from the startpoint
! nax is number of axis, axarr are description of the axis
! axarr are axis records
! tmpline is a line record ... not needed ... ??
    implicit none

```

```

integer nax,axactive,ieq
type(meq_setup), pointer :: meqrec
type(map_line), dimension(2) :: tmpline
type(map_axis), dimension(nax) :: axarr
type(gtp_equilibrium_data), pointer :: ceq

```

5.5 Storing a calculated point on the line

```

subroutine map_store(mapline,axarr,nax,saveceq)
! store a calculated equilibrium
  implicit none
  integer nax
  type(map_line), pointer :: mapline
  type(map_axis), dimension(nax) :: axarr
  type(map_ceqresults), pointer :: saveceq

```

5.6 Storing the last equilibrium along a line

```

subroutine map_lineend(mapline,value,ceq)
! terminates gracefully a line at an axis limit or an error.
! maptop probably not needed except for testing
  implicit none
  integer mode
  type(map_line), pointer :: mapline
  type(gtp_equilibrium_data), pointer :: ceq
  double precision value

```

5.7 Changing the axis for next equilibrium calculation

During mapping it may sometimes be necessary to change the axis variable used for moving along the line.

```

subroutine map_changeaxis(mapline,nyax,oldax,nax,axarr,axval,bytax,ceq)
! changes the axis with active condition to nyax
  type(map_line), pointer :: mapline
  type(gtp_equilibrium_data), pointer :: ceq
  type(map_axis), dimension(nax) :: axarr
  logical bytax
! nax is number of axis, nyax is new axis with active condition
  integer nyax,nax,oldax
! the value to set as condition on new axis

```

```

    double precision axval
    subroutine map_force_changeaxis(maptop,mapline,meqrec,nax,axarr,ceq)
! force change of axis with active condition. Works only with 2 axis.
! (and for tie-line not in plane ??). Similar to map_changeaxis ...
    implicit none
! number of axis, also in maptop record
    integer nax
    type(map_node), pointer :: maptop
    type(map_line), pointer :: mapline
    type(meq_setup) :: meqrec
    type(gtp_equilibrium_data), pointer :: ceq
    type(map_axis), dimension(*) :: axarr

```

5.8 Calculating the next equilibrium along a line

This is the normal routine to calculate a new equilibrium along a line.

```

    subroutine map_step(maptop,mapline,meqrec,phr,axvalok,nax,axarr,ceq)
! used also for map as mapping is stepping in one axis with phase fix condition
! calculate the next equilibrium along a line. New phases can appear.
! axis with active condition can change and the direction.
    implicit none
! number of axis, redundant as also in maptop record
    integer nax
    type(map_node), pointer :: maptop
    type(map_line), pointer :: mapline
    type(meq_setup) :: meqrec
! phr is included just for debugging to see phase amounts
    type(meq_phase), dimension(*), target :: phr
    type(gtp_equilibrium_data), pointer :: ceq
    type(map_axis), dimension(*) :: axarr
    double precision axvalok

```

5.9 Calculating a node point

This is used to calculate a node point when the set of phases are changing.

```

    subroutine map_calcnode(irem,iadd,maptop,mapline,meqrec,axarr,ceq)
! we have found a change in the set of stable phases. check if this node
! already been found and if so eliminate a line record. Otherwise
! create a new node record with line records and continue mapping one

```

```

! of these.
    implicit none
    integer irem,iadd
! am am wondering if pointer is necessary??
    type(map_node), pointer :: maptop
    type(map_line), pointer :: mapline
! Note changes in meqrec is local, not copied to mapline%meqrec!!!
    type(meq_setup) :: meqrec
    type(map_axis), dimension(*) :: axarr
    type(gtp_equilibrium_data), pointer :: ceq

```

5.10 Creates a node with several lines

After calculating the node point this routine determines the sxits for new lines.

```

    subroutine map_newnode(mapline,meqrec,maptop,axval,lastax,axarr,phfix,ceq)
! must be partially THREADPROTECTED
! first check if a node with this equilibrium already exists
! if not add a new node with appropriate lineheads and arrange all links
! Take care it tie-lines in the plane all lines do not have to be calculated
! NOTE: meqrec not the same as mapline%meqrec !! ??
    type(map_node), pointer :: maptop
    type(meq_setup) :: meqrec
    type(map_line), pointer :: mapline,nodexit
    type(map_axis), dimension(*) :: axarr
    type(gtp_equilibrium_data), pointer :: ceq
    integer phfix,lastax
! axval is axis value which was attempted to calculate,
    double precision axval

```

5.11 Reserves a record to save a calculated equilibrium

```

    subroutine reserve_saveceq(location,saveceq)
! must be THREADPROTECTED
! reserves a ceq record in saveceq
    implicit none
    integer location
    type(map_ceqresults), pointer :: saveceq

```

5.12 Finds a line to be calculated

Searches all node points for lines to be calculated.

```
subroutine map_findline(maptop,axarr,mapfix,mapline)
! must be THREADPROTECTED
! Find a map_line record to be calculated from maptop
! already been found and if so eliminate a line record. Otherwise
  type(map_node), pointer :: maptop
  type(map_line), pointer :: mapline
  type(map_axis), dimension(*) :: axarr
  type(map_fixph), pointer :: mapfix
```

5.13 Creates records for saving equilibria

```
subroutine create_saveceq(ceqres,size)
! creates an array of equilibrium records to save calculated lines for step
! and map
  type(map_ceqresults), pointer :: ceqres
  integer size
```

5.14 Checks if the diagram has tie-lines in the plane

```
integer function tieline_inplane(nax,axarr,ceq)
! returns -1 if tielines are not in the plane (isopleth)
!           0 for step calculations (nax=1)
!           1 if tielines in the plane (binary T-X, ternary isopleths
!           set if more than one extensive variable is not axis variables
  integer nax
  type(map_axis), dimension(nax) :: axarr
  type(gtp_equilibrium_data), pointer :: ceq
```

5.15 Checks if the node point is an invariant equilibrium

A nodepoint from an invariant equilibria have more lines that exit than normal nodes.

```
integer function invariant_equilibrium(lines,mapnode)
! Only called for tie-lines not in plane. If tie-lines in plane then all
! nodes are invariants.
  integer lines
  type(map_node), pointer :: mapnode
```


5.16 Tries to handel problems when mapping

This tries various way to handle problems during STEP and MAP.

```
subroutine map_problems(maptop,mapline,axarr,xxx,typ)
! jump here for different problems
  integer typ
  type(map_node), pointer :: maptop
  type(map_line), pointer :: mapline
  type(map_axis), dimension(*) :: axarr
  double precision xxx,yyy
```

5.17 Using smaller steps when there are convergence problems

This subroutine is used when there is a convergence problem calculating an equilibrium.

```
subroutine map_halfstep(halfstep,axvalok,mapline,axarr,ceq)
! Used when an error calculating a normal step or a node point
! take back the last sucessfully calculated axis value and take smaller step
! possibly one should also restore the ceq record.
  implicit none
  integer halfstep
  double precision axvalok
  TYPE(gtp_equilibrium_data), pointer :: ceq
  TYPE(map_line), pointer :: mapline
  type(map_axis), dimension(*) :: axarr
```

5.18 Lists stored equilibria

```
subroutine list_stored_equilibria(kou,axarr,maptop)
! list all nodes and lines from step/map
! maybe allow some delete/amend later ...
  integer kou
  type(map_node), pointer :: maptop
  type(map_axis), dimension(*) :: axarr
```

5.19 Plots calculated diagrams

```
subroutine ocplot2(ndx,pltax,filename,maptop,axarr,graphopt,pform,ceq)
```

```

! ndx is number of plot axis, pltax is text with plotaxis variables
! filename is intermediary file (maybe not needed)
! maptop is map_node record with all results
! pform is type of output (screen or postscript or jpeg)
    implicit none
    integer ndx
    character pltax(*)*(*),filename*(*),pform*(*)
    type(map_axis), dimension(*) :: axarr
    type(map_node), pointer :: maptop
    type(graphics_options) :: graphopt

```

5.20 Calculates equilibria for one phase at a time

Needed to calculate for example Gibbs energy curves.

```

subroutine step_separate(maptop,noofaxis,axarr,starteq)
! calculates for each phase separately along an axis (like G curves)
! There can not be any changes of the stable phase ...
    implicit none
    integer noofaxis
    type(map_axis), dimension(noofaxis) :: axarr
    TYPE(gtp_equilibrium_data), pointer :: starteq
    TYPE(map_node), pointer :: maptop

```

6 Summary

That is all!

References

- [84Mor] J.E. Morall, Scripta Metallurgica, 18, (1984), 407
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- [15Sun2] B Sundman, X-G Lu, H Ohtani, Comp Mat. Sci **101**(2015) 127