

# XPP Commands

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## ODE File Format

```
# comment line - name of file, etc
#include filename
d<name>/dt=<formula>
<name>'=<formula>
<name>(t)=<formula>
volt <name>=<formula>
<name>(t+1)=<formula>
x[n1..n2]' = ...[j] [j-1] ... <-- Arrays
%[i1..i2]
u[j]'=...
v[j]'=...
%
markov <name> <nstates>
    {t01} {t02} ... {t0k-1}
    {t10} ...
    ...
    {tk-1,0} ... {tk-1 k-1}

aux <name>=<formula>
!<name>=<formula> <-- parameters defined as formulae
<name>=<formula>
parameter <name1>=<value1>,<name2>=<value2>, ...
wiener <name1>, <name2>, ...
number <name1>=<value1>,<name2>=<value2>, ...
<name>(<x1>,<x2>,...,<xn>)=<formula>
table <name> <filename>
table <name> % <npts> <xlo> <xhi> <function(t)>
global sign {condition} {name1=form1;...}
init <name>=<value>,...
<name>(0)=<value> or <expr> <-- delay initial conditions
bdry <expression>
0= <expression> <--- For DAEs
solv <name>=<expression> <----- For DAEs
special <name>=conv(type,npts,ncon,wgt,rootname)
        fconv(type,npts,ncon,wgt,rootname,root2,function)
        sparse(npts,ncon,wgt,index,rootname)
        fsparse(npts,ncon,wgt,index,rootname,root2,function)
# comments
@ <name>=<value>, ...
set <name> {x1=z1,x2=z2,...}
only <name1>,<name2>,...
options <filename>
" {z=3,b=3,...} Some nice text <--- Active comments
done
```

### INTEGRAL EQUATIONS

The general integral equation

$$u(t) = f(t) + \int_0^t K(t,s,u(s))ds$$

becomes

$$u = f(t) + \int K(t, t', u) dt'$$

The convolution equation:

$$v(t) = \exp(-t) + \int_0^t e^{-(t-s)} v(s) ds$$

would be written as:

$$v(t) = \exp(-t) + \int \exp(-t^2) v dt$$

If one wants to solve, say,

$$u(t) = \exp(-t) + \int_0^t (t-t')^{-mu} K(t, t', u(t')) dt'$$

the form is:

$$u(t) = \exp(-t) + \int [\mu] K(t, t', u) dt'$$

and for convolutions, use the form:

$$u(t) = \exp(-t) + \int [\mu] w(t) u dt$$

## NETWORKS

`special zip=conv(type,npts,ncon,wgt,root)`

where `root` is the name of a variable and `wgt` is a table, produces an array `zip` with `npts`:

$$zip[i] = \sum_{j=-ncon}^{ncon} wgt[j + ncon] root[i + j]$$

The `sparse` network has the syntax:

`special zip=sparse(npts,ncon,wgt,index,root)`

where `wgt` and `index` are tables with at least `npts * ncon` entries. The array `index` returns the indices of the offsets to which to connect and the array `wgt` is the coupling strength. The return is

$$zip[i] = \sum_{j=0; j < ncon} w[i * ncon + j] * root[k] \\ k = index[i * ncon + j]$$

The other two types of networks allow more complicated interactions:

`special zip=fconv(type,npts,ncon,wgt,root1,root2,f)`

evaluates as

$$zip[i] = \sum_{j=-ncon; j < ncon} wgt[ncon + j] * f(root1[i + j], root2[i])$$

and

`special zip=fsparse(npts,ncon,wgt,index,root1,root2,f)`

evaluates as

$$zip[i] = \sum_{j=0; j < ncon} wgt[ncon * i + j] * f(root1[k], root2[i]) \\ k = index[i * ncon + j]$$

Matrix multiplication is also possible:

```
special k=mmult(n,m,w,u)
```

returns a vector **k** of length **m** defined as

```
k(j)=sum(i=0;i<n)w(i+nj)u(i)
```

while

```
special k=fmmult(n,m,w,u,v,f)
```

returns

```
k(j)=sum(i=0;i<n)w(i+nj)f(u(i),v(j))
```

**OPTIONS** The format for changing the options is:

```
@ name1=value1, name2=value2, ...
```

where **name** is one of the following and **value** is either an integer, floating point, or string. (All names can be upper or lower case). The first four options *can only be set outside the program*. They are:

- **MAXSTOR=integer** sets the total number of time steps that will be kept in memory. The default is 5000. If you want to perform very long integrations change this to some large number.
- **BACK= {Black,White}** sets the background to black or white.
- **SMALLFONT=fontname** where **fontname** is some font available to your X-server. This sets the “small” font which is used in the Data Browser and in some other windows.
- **BIGFONT=fontname** sets the font for all the menus and popups.
- **SMC={0,...,10}** sets the stable manifold color
- **UMC={0,...,10}** sets the unstable manifold color
- **XNC={0,...,10}** sets the X-nullcline color
- **YNC={0,...,10}** sets the Y-nullcline color
- **SEC,UEC,SPC,UPC=color** set the colors for the screen on AUTO for stable eq, unstable eq, stable per, unstable per.
- **OUTPUT=filename** sets the filename to which you want to write for “silent” integration. The default is “output.dat”.
- **DWCOLOR,MWCOLOR,BACKCOLOR,FORECOLOR=HEXCOLOR**. Set drawing window,mainwindow,background color,foreground color as a hexadecimal, e.g. FF0000 is pure red, 090909 is medium grey, 000000 is black, FFFFFFFF is white.
- **GRADS={0,1}** turn off or on gradients in the XPP buttons  $j/LI_i$
- **HEIGHT=pixels,WIDTH=pixels**, sets the height and width of the main window initially
- **RUNNOW=1** run the simulation as soon as the windows are up (Don’t wait for Init conds Go, e.g.
- **BUT=name:kbs** defines button on the top of the main window. You can define up to 20 such buttons. They will appear across the top when you press them, they will execute an XPP keyboard shortcut. For example, **BUT=Fit:wf** will create a button labeled **Fit** and when you press it, it will be as if you had clicked **Window/zoom Fit** Most of the menu items are available.

The remaining options can be set from within the program. They are

- **LT=int** sets the linetype. It should be less than 2 and greater than -6.
- **SEED=int** sets the random number generator seed.
- **XP=name** sets the name of the variable to plot on the x-axis. The default is **T**, the time-variable.
- **YP=name** sets the name of the variable on the y-axis.
- **ZP=name** sets the name of the variable on the z-axis (if the plot is 3D.)
- **COLORMAP=0,1,2,3,4,5** sets the colormap
- **NPLOT=int** tells XPP how many plots will be in the opening screen.

- `XP2=name,YP2=name,ZP2=name` tells XPP the variables on the axes of the second curve; `XP8` etc are for the 8th plot. Up to 8 total plots can be specified on opening. They will be given different colors.
- `AXES={2,3}` determine whether a 2D or 3D plot will be displayed.
- `TOTAL=value` sets the total amount of time to integrate the equations (default is 20).
- `DT=value` sets the time step for the integrator (default is 0.05).
- `NJMP=integer` tells XPP how frequently to output the solution to the ODE. The default is 1, which means at each integration step.
- `T0=value` sets the starting time (default is 0).
- `TRANS=value` tells XPP to integrate until `T=TRANS` and then start plotting solutions (default is 0.)
- `NMESH=integer` sets the mesh size for computing nullclines (default is 40).
- `METH={ discrete,euler,modeuler,rungekutta,adams,gear,volterra, backeul, qualrk,stiff,cvode,5dp,` sets the integration method (default is Runge-Kutta.)
- `DTMIN=value` sets the minimum allowable timestep for the Gear integrator.
- `DTMAX=value` sets the maximum allowable timestep for the Gear integrator
- `VMAXPTS=value` sets the number of points maintained in for the Volterra integral solver. The default is 4000.
- `{ JAC_EPS=value, NEWT_TOL=value, NEWT_ITER=value}` set parameters for the root finders.
- `ATOLER=value` sets the absolute tolerance for CVODE.
- `TOLER=value` sets the error tolerance for the Gear, adaptive RK, and stiff integrators. It is the relative tolerance for CVODE.
- `BOUND=value` sets the maximum bound any plotted variable can reach in magnitude. If any plottable quantity exceeds this, the integrator will halt with a warning. The program will not stop however (default is 100.)
- `DELAY=value` sets the maximum delay allowed in the integration (default is 0.)
- `BANDUP=int, BANDLO=int` bandwidths for the Jacobian computation.
- `PHI=value,THETA=value` set the angles for the three-dimensional plots.
- `XLO=value,YLO=value,XHI=value,YHI=value` set the limits for two-dimensional plots (defaults are 0,-2,20,2 respectively.) Note that for three-dimensional plots, the plot is scaled to a cube with vertices that are  $\pm 1$  and this cube is rotated and projected onto the plane so setting these to  $\pm 2$  works well for 3D plots.
- `XMAX=value, XMIN=value, YMAX=value, YMIN=value, ZMAX=value, ZMIN=value` set the scaling for three-d plots.
- `POIMAP={ section,maxmin, period}` sets up a Poincare map for either sections of a variable, the extrema, or period.
- `POIVAR=name` sets the variable name whose section you are interested in finding.
- `POIPLN=value` is the value of the section; it is a floating point.
- `POISGN={ 1, -1, 0 }` determines the direction of the section.
- `POISTOP=1` means to stop the integration when the section is reached.
- `RANGE=1` means that you want to run a range integration (in batch mode).
- `RANGEOVER=name, RANGESTEP=number, RANGELOW=number, RANGEHIGH=number, RANGERESET=Yes,No, RANGEOLDIC=Yes,No` all correspond to the entries in the range integration option.
- `TOR_PER=value`, defined the period for a toroidal phasespace and tellx XPP that there will be some variables on the circle.
- `FOLD=name`, tells XPP that the variable `name` is to be considered modulo the period. You can repeat this for many variables.
- **AUTO-stuff.** The following AUTO-specific variables can also be set: `NTST, NMAX, NPR, DSMIN, DSMAX, DS, PARMIN, PARMAX, NORMMIN, NORMMAX, AUTOXMIN, AUTOXMAX, AUTOYMIN, AUTOYMAX, AUTOVAR`. The last is the variable to plot on the y-axis. The x-axis variable is always the first parameter in the ODE file unless you change it within AUTO.
- `DLL_LIB=file` Dynamically linked library
- `DLL_FUN=name` Dynamically linked function.

**COLOR MEANING** 0-Foreground color; 1-Red; 2-Red Orange; 3-Orange; 4-Yellow Orange; 5-Yellow; 6-Yellow Green; 7-Green; 8-Blue Green; 9-Blue; 10-Purple.

**KEYWORDS** You should be aware of the following keywords that should not be used in your ODE files for anything other than their meaning here.

```
sin cos tan atan atan2 sinh cosh tanh
exp delay ln log log10 t pi if then else
asin acos heav sign ceil flr ran abs del\_shift
max min normal besselj bessely besseli erf erfc poisson
lgamma arg1 ... arg9 @ $ + - / * ^ ** shift
| > < == >= <= != not \# int sum of i'
```

These are mainly self-explanatory. The nonobvious ones are:

- **heav(arg1)** the step function, zero if **arg1**<0 and 1 otherwise.
- **sign(arg)** which is the sign of the argument (zero has sign 0)
- **ran(arg)** produces a uniformly distributed random number between 0 and **arg**.
- **besselj**, **bessely**, **besseli** take two arguments,  $n, x$  and return  $J_n(x)$ ,  $Y_n(x)$ ,  $I_n(x)$  the Bessel functions.
- **erf(x)**, **erfc(x)** are the error function and the complementary function.
- **lgamma(x)** is the log of the gamma function.
- **normal(arg1,arg2)** produces a normally distributed random number with mean **arg1** and variance **arg2**.
- **poisson(arg)** produces the number of events from a Poisson process with parameter **arg**. It is a random number.
- **max(arg1,arg2)** produces the maximum of the two arguments and **min** is the minimum of them.
- **if(<exp1>)then(<exp2>)else(<exp3>)** evaluates **<exp1>**. If it is nonzero it evaluates to **<exp2>** otherwise it is **<exp3>**. E.g. **if(x>1)then(ln(x))else(x-1)** will lead to **ln(2)** if **x=2** and **-1** if **x=0**.
- **delay(<var>,<exp>)** returns variable **<var>** delayed by the result of evaluating **<exp>**. In order to use the delay you must inform the program of the maximal possible delay so it can allocate storage.
- **del\_shift(<var>,<shift>,<delay>)**. This operator combines the **delay** and the **shift** operators and returns the value of the variable **<var>** shifted by **<shift>** at the delayed time given by **<delay>**.
- **ceil(arg)**, **flr(arg)** are the integer parts of **<arg>** returning the smallest integer greater than and the largest integer less than **<arg>**.
- **t** is the current time in the integration of the differential equation.
- **pi** is  $\pi$ .
- **arg1**, ..., **arg9** are the formal arguments for functions
- **int**, **#** concern Volterra equations.
- **shift(<var>,<exp>)** This operator evaluates the expression **<exp>** converts it to an integer and then uses this to indirectly address a variable whose address is that of **<var>** plus the integer value of the expression. This is a way to imitate arrays in XPP. For example if you defined the sequence of 5 variables, **u0,u1,u2,u3,u4** one right after another, then **shift(u0,2)** would return the value of **u2**.
- **sum(<ex1>,<ex2>)of(<ex3>)** is a way of summing up things. The expressions **<ex1>**, **<ex2>** are evaluated and their integer parts are used as the lower and upper limits of the sum. The index of the sum is **i'** so that you cannot have double sums since there is only one index. **<ex3>** is the expression to be summed and will generally involve **i'**. For example **sum(1,10)of(i')** will be evaluated to 55. Another example combines the sum with the shift operator. **sum(0,4)of(shift(u0,i'))** will sum up **u0** and the next four variables that were defined after it.

Call **xppaut -help** to get the command line arguments. There are many.

Add a **.xpprc** file to set your favorite options, e.g

```
@ bell=0,grads=0,dwcolor=eeddff
@ bigfont=lucidasanstypewriter-bold-14
```