



A Publication for ANSYS Users

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# The Focus



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## Tools of Post Processing: Listing

By [Eric Miller](#), PADT

Post processing is an oft-maligned part of the analysis process. Real analysts focus on meshing and running their models. But the fact of the matter is that getting *meaningful answers* is the whole point of doing a simulation, so spending a little bit of time to learn how to produce more concise and digestible results is a good investment. To that end, we will be covering a couple of key ways to produce higher quality results using POST1 and POST26, including plotting, animation, sections, and for this first article, listing.



### Listing Results

Because of ANSYS' long history, it has a rich listing capability. Back in the old days, listing was the primary way to extract information. But even now, it is still an important way to nail down facts about your results. By default, the formatting is not very clean and it is easy to get buried in details. Learning a few commands can turn that pile of numbers into valuable tables and eliminate considerable manual cleanup in external programs.

The first thing to know about ANSYS listings is that they are set up for sending data directly to large volume line printers used 10 or 15 years ago. That's why that silly header shows up every 25 lines. It's also the reason why the information is presented in unformatted Courier1 font without tabs, separating lines, bolding, or italics. Old printers just could not produce that type of formatting. So don't try to dress up listings by changing font or style unless you insert tabs into your data first.

The most important formatting command is (*surprise!*), `/format`:

`/FORMAT, NDIGIT , Ftype, NWIDTH, DSIGNF, LINE, CHAR>`

- *NDIGIT*: Number of digits in first column (usually entity number)
- *Ftype*: FORTRAN format type to use: G, F, or E
- *NWIDTH*: Total width of number
- *DSIGNF*: Significant figures
- *LINE*: Number of lines per page (defaults to `/page`)
- *CHAR*: Number of characters per line before wrap (defaults to `/page`)

Figure 1 shows listings with the default formatting and with controls placed on spacing. The listing wraps and is hard to read with the defaults (formatted for a 132 column line-printer.) You could also use the *DSIGNF* argument if you need more precision on a number.



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Default /format (Windows)											
PRINT 5    NODEL SOLUTION PER NODE											
NODE	UX	UY	UZ	UXY	UYZ	UXZ	S1	S2	S3	SHEAR	SDIV
1	5.8829	295.41	-15.332	11.429	-29.218	8.1254E-13	291.31	5.2263	-16.455	315.65	384.43
2	-55.683	399.19	-29.129	-6.9541	-21.458	-54.454	488.32	2.8228	-61.695	461.82	451.35
3	4.5851	265.62	-1.8656	16.483	38.882	1.3833	269.98	5.1495	-11.528	283.25	214.28
4	-15.448	218.31	-168.85	-15.888	18.558	5.1198	213.14	-16.355	-168.95	452.69	583.49

/format,8,f,9,1,100,240											
PRINT 5    NODEL SOLUTION PER NODE											
NODE	UX	UY	UZ	UXY	UYZ	UXZ	S1	S2	S3	SHEAR	SDIV
1	5.1	295.4	-15.3	11.4	-29.2	8.1	291.2	5.2	-16.5	315.6	384.4
2	-55.6	399.1	-29.1	-6.9	-21.4	-54.5	488.3	2.8	-61.1	461.8	451.3
3	4.6	265.6	-1.9	16.4	38.8	1.2	269.9	5.1	-11.5	283.2	214.2
4	-15.4	218.3	-168.8	-15.8	18.5	5.1	213.1	-16.2	-168.8	452.1	583.5
5	-15.1	-345.3	51.4	-2.3	-6.8	-82.9	313.4	-69.6	-345.2	456.6	398.5
6	-4.8	-369.8	238.2	-2.6	-9.2	54.5	233.8	-11.5	-369.2	688.2	522.5
7	11.5	614.8	-85.8	5.4	1.8	11.5	614.9	14.4	-89.8	183.9	658.4

Figure 1.

The next area that needs user-tweaking is cleaning up the header. The /HEADER command contains a series of on/off switches that can be used to simplify and reduce the amount of information printed at the top of data listings. It is also very useful when you want to dump a listing to a file that will be read by another program. This is because you can turn off that pesky header information that is such a pain to skip over. The command is fairly simple:

/HEADER, *Header*, *Stitle*, *Ldstmp*, *Notes*, *Colhed*, *Minmax*

- *Header*: Standard ANSYS header (batch only, always off in interactive mode)
- *Stitle*: Subtitles
- *Ldstmp*: Load step information for listed data
- *Notes*: Information on a particular table listing
- *Colhed*: Column headings on table listings
- *Minmax*: Min/Max information at the end of tables

Figure 2 shows the default header with all that extra junk and spaces, as well as two alternative listings that are much more readable.



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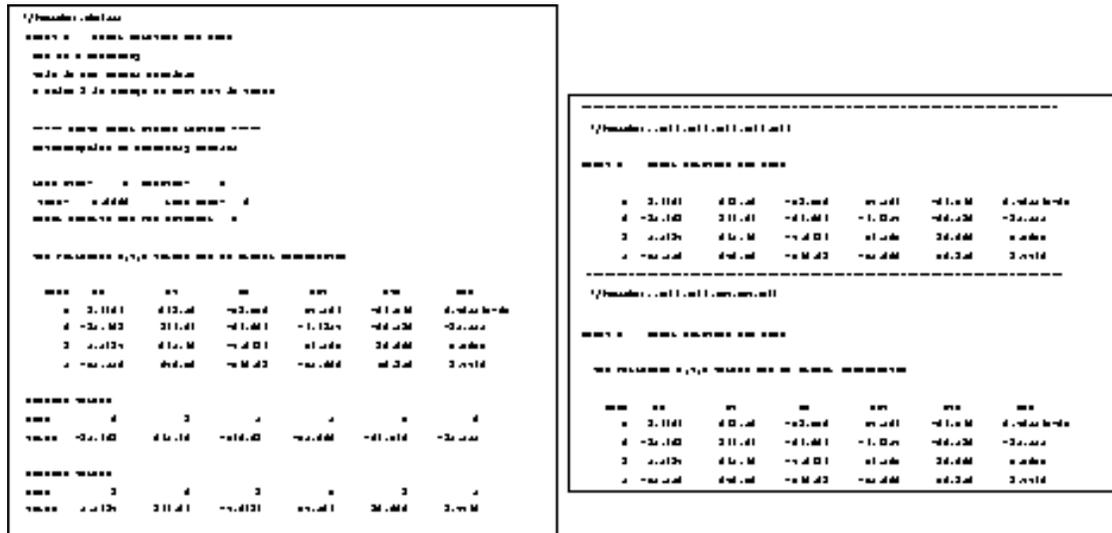


Figure 2.

However, this does not solve the most annoying aspect of saving listings when running in batch mode to a file – a header every 24 lines. You control this with the `/PAGE` command:

`/PAGE, ILINE, ICHAR, BLINE, BCHAR`

- *ILINE, ICHAR*: no longer used
- *BLINE*: Number of lines per page for batch, /output or interactive listing
- *BCHAR*: Number of characters per line for batch, /output or interactive listing

I set *BLINE* to 9999 and get only one header.

But what if ANSYS does not list the data you want, or does not list it in a way you find useful. Since you can extract information from the database with `*GET`'s and `*VGET`'s you can pretty much format information any way you want. The most common is to do a `*VGET` followed by `*CFOPEN`, `*VWRITE`'s, and `*CFCLOSE`. `*VWRITE` allows you to enter FORTRAN and C format statements. There is also an `*MWRITE` command for outputting matrices. With these, output to programs like Microsoft Excel is a breeze. Figure 3 shows a very simple macro that outputs node numbers and X locations for 10 nodes using various formats.

```
finish
/clear
/prep7
n,1
n,10,1
fill,1,10

*dim,nnx,,10,2
*vfill,nnx(1,1),ramp,1,1
*vget,nnx(1,2),node,1,loc,x

*cfopen,vwrt,txt
```



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```

*vwrite,
----- Comma Separated 'C':
*vwrite,
(" ")
*vwrite,nnx(1,1),nnx(1,2)
%8d,%16.9g

*vwrite,
(" ")
*vwrite,
----- Comma Separated FORTRAN:
*vwrite,
(" ")
*vwrite,nnx(1,1),nnx(1,2)
(F8.0,"",G16.9)

*vwrite,
(" ")
*vwrite,
----- Text in the line:
*vwrite,
(" ")
*vwrite,nnx(1,1),nnx(1,2)
Node: %8d X-Position: %16.9g

*vwrite,
(" ")
*vwrite,
----- FORTRAN Again
*vwrite,
(" ")
*vwrite,nnx(1,1),nnx(1,2)
(" -- ",F8.0,"": ",E16.9)
*cfclose
*ulist,vwrt,txt

```

Figure 3. Example macro.

```

----- Comma Separated 'C':

1, 0.000000000
2, 0.250000000
3, 0.500000000
4, 0.750000000
5, 1.000000000

```



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----- Comma Separated FORTRAN:

```
1., 0.000000000
2., 0.250000000
3., 0.500000000
4., 0.750000000
5., 1.000000000
```

----- Text in the line:

```
Node:    1 X-Position:  0.000000000
Node:    2 X-Position:  0.250000000
Node:    3 X-Position:  0.500000000
Node:    4 X-Position:  0.750000000
Node:    5 X-Position:  1.000000000
```

----- FORTRAN Again

```
--      1: 0.000000000E+00
--      2: 0.250000000E+00
--      3: 0.500000000E+00
--      4: 0.750000000E+00
--      5: 0.100000000E+01
```

Figure 4. Output in vwrt.txt.

Take note that the first example is comma separated. A \*VWRITE or \*MWRITE with commas has proven to be the best way to output to Microsoft Excel. I prefer the C formatting because it is more concise and allows for integers. For various reasons, the FORTRAN (I) format does not work well on some platforms.

However, if that is not enough control and you want to include fancy formatting and your own text headers, the best solution is to write an external command (~command) or use the built in Tcl/Tk interpreter. These options allow you to extract information with \*VGET's, and then output them with the full formatting provided by a modern programming language.

In short, you can pretty much format output any way you want with a little bit of set up, especially with the \*VWRITE command. I have found this especially useful when I am automating report generation or when I'm doing a parametric study. Scripting the format of the listing allows me to have consistent and readable output. Some say the days of six-foot high stacks of result printouts are gone, but that does not mean that you cannot find value in outputting values.

---

1 Courier belongs to a family of fonts called "monotype" or "fixed width". What this means is that each character is exactly the same width. Most fonts are "proportional", which means they



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have characters that vary in width for aesthetic reasons, which makes formatting without tabs a real pain. To illustrate, here are two examples of some lines of text The first pair is in Times New Roman (a proportional font) and the second is in Courier. Compare the ends of the lines.

This line has exactly forty characters!!  
Hello, I am forty characters long. Even?

This line has exactly forty characters!!  
Hello, I am forty characters long. Even?



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# User Programmable Features

by [Eric Miller](#), PADT, and [Rod Scholl](#), PADT

## UPFs: What Are They?

One of ANSYS' strengths is that its toolbox will handle anything from a textbook plate with hole, all the way up to allowing you to recompile the code with FORTRAN scripts that you develop yourself. The User Programmable Feature (UPF) capability allows users to create their own subroutines and link them into ANSYS.

## What Are UPFs Used For?

There are a variety of common uses for UPFs:

- Material models
- User-developed elements
- User commands
- Element orientation
- User beam definition
- Heat generation
- Defining real constants "on the fly"



## How Hard Is It to Implement a UPF?

Implementing a UPF is definitely more work than simply programming the action using APDL. However, if you have existing FORTRAN code you want to include or run externally, or have any of the above-mentioned challenges, then APDL will not suffice.

To learn for yourself, you will want to surround your self with resources:

- This article!



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- *ANSYS Programmer's Guide: User Programmable Features*
  - A **must-have!!!!**
  - All the needed subroutines are documented here
- The hundreds of documented template scripts found within the installation directory in /custom/user
- [XANSYS](#). Submit your questions to the XANSYS community.
- PADT. Submit your questions to us.

## What is the Process?

1. Copy a sample routine from the ANSYS directories
2. Modify it to do what you want
3. Use the ANSYS batch compile script to compile and link
4. Locate the new executable for use
5. In your model, activate user features by applicable method
  - a. Define user element
  - b. Define user material model
  - c. Turn on user feature in element real constants
  - d. Define user command with /UCMD
  - e. Setup in USRCAL command
6. Debug



## Do I have to use FORTRAN?

- All of these routines are called as FORTRAN and use FORTRAN calls to utility functions.
- You could use FORTRAN wrappers to call C routines, but that is usually not worth the effort.



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- ANSYS provides you with a compile script to help with the process.
- Don't freak out. FORTRAN is very easy to use, and creating UPFs and recompiling ANSYS is actually very easy.

## You Want to Learn by Example!?!?

So before opening the *ANSYS Programmer's Guide*, you want to take a look at what you're getting into? Well then, buckle up...

### UPFs Example1, USER02.F

- USER02.F contains code to offset nodes by supplied x,y,z values
- Good example of structure of routines
- Shows how other routines are called

**Typical Header Information.** Includes source control data (\*deck), copyright information, confidentiality, and pointer to more information. Also function statement.

```
*deck,user02      user
ANSYS,INC
  function user02 (initin,dpin,ch4in,ch8in)
c *** primary function:  user routine number 02
c   This demonstration offsets selected nodes with the command:
c   usr2,dx,dy,dz

c   *** ansys (r) copyright(c) 2000
c   *** ansys, inc.
c *** Notice - This file contains ANSYS Confidential information

c /*****\
c | see user01 for additional information on user routines   |
c \*****/
<snip>
```

**Arguments Section.** Describes input and output arguments, what type they are, and what they are used for. Also, additional examples and help are put in this section. Externally passed and local variables are also defined here. Finally, the standard include statements are placed here.



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<snip>

c input arguments:

c variable (typ,siz,intent) description

c intin (int,ar(12),in) - integers from command line

c dpin (dp,ar(12),in) - double precision from cmnd line

c ch4in (ch\*4,ar(12),in) - upper case 4 characters from cmnd

c ch8in (ch\*8,ar(12),in) - as input 8 characters from cmnd

c output arguments: none

c user02 (int,sc,out) - result code (should be zero)

c (which is ignored for now)

c \*\*\*\*\*

c intin(2), dpin(2), ch4in(2), & ch8in are all representations

c of the contents of the second field on the command line

c (the field after the command label)

c \*\*\*\*\*

external wrinqr,ndinqr,ndgxyz,ndpxyz,erhandler

integer wrinqr,ndinqr,ndgxyz

integer usr02,intin(12),iott,maxnp,i,ksel

double precision dpin(12),xyz(3)

character\*4 ch4in(12)

character\*8 ch8in(12)

#include "ansysdef.inc"

<snip>

**Calculations.** Use ndinqr() to get max node number. This code snippet loops on all possible node numbers, checks to see if the node is selected, and gets its *x*, *y*, and *z* values. Then it does the offset and updates the *x,y,z* position with ndpxyz().

<snip>

maxnp = ndinqr (0,DB\_MAXDEFINED)

do i = 1,maxnp

ksel = ndgxyz (i,xyz(i))

if (ksel .eq. 1) then

xyz(1) = xyz(1) + dpin(2)



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```
xyz(2) = xyz(2) + dpin(3)
xyz(3) = xyz(3) + dpin(4)
call ndpxyz (i,xyz(i))
endif
enddo
<snip>
```

**Clean Up and Exit.** This code snippet is used to figure out the standard output unit, and then write a message to it. It will also write a message to a GUI pop-up box. It will then set the return value and leave.

```
<snip>
c ***** write to output file *****
  iott = wrinqr(WR_OUTPUT)
  write (iott,2000)
  2000 format (' NODE OFFSET COMPLETE '/')

c ***** write to GUI window *****
  call erhandler ('user02',3000,
x   2,'NODE OFFSET COMPLETE',0.0d0,' ')

c ***** required return value *****
  user02 = 0

  return
end
<snip>
```

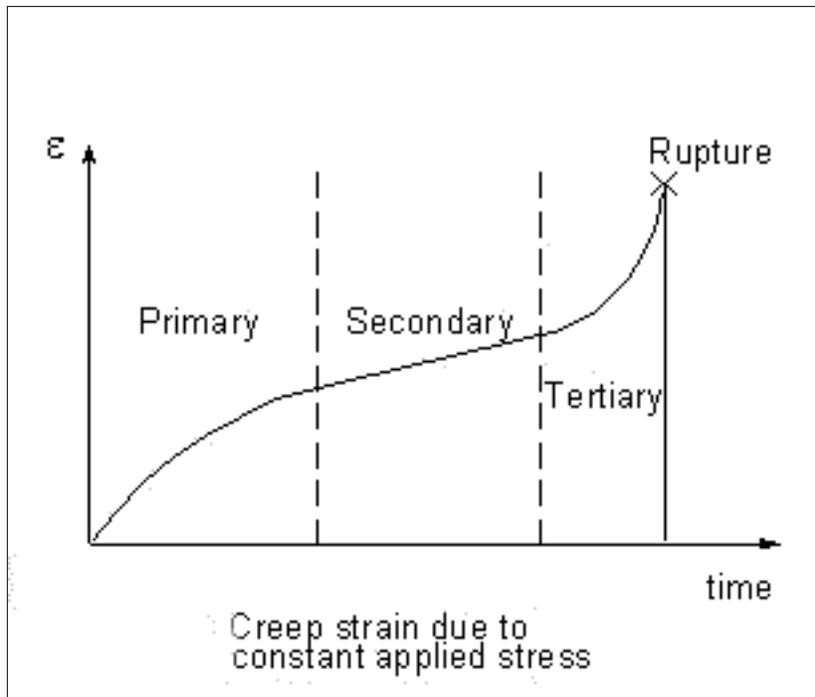
## UPFs: Example 2, USERCR.F

Creep strain due to constant applied stress:

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```
subroutine usercr (elem,intpt,mat,ncomp,kfirst,kfsteq,e,posn,d,
  x proptb,timval,timinc,tem,dtem,tofst,fluen,dfluen,epel,
  x epcrp,statev,usvr,delcr)
```

```
c
```

```
#include "impcom.inc"
  external erhandler
```

```
c
```

```
c user-defined fortran parameters
```

```
c --- size of usvr data
  integer nuval,nintp
  parameter (nuval=1,nintp=1)
```

```
c
```

```
c external subroutines and functions
```

```
  external egen
  double precision egen
  external vapb1,vamb1,vmult
```

```
c
```

```
c integer variables
```

```
  integer elem,intpt,mat,ncomp,kfirst,kfsteq
```

```
c
```

```
c double precision variables
```

```
  double precision
```



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```

x e, posn, d(ncomp, ncomp), proptb(72), timval, timinc, tem, dtem,
x tofst, fluen, dfluen, epel(ncomp), epcrp(ncomp),
x statev(ncomp*5+2), usvr(nuval, nintp), delcr, temabs, con,
x del(6), epet, sigen, eptot, ept(6),
c -PADT--- PADT added variables
x qovrk, c1, alpha, dbln, crprt, vrbflg,
x aaaa, bbbb, cccc, dddd
<snip>
<snip>
c
c --- initial checks
c --- author should remove the warning below when making changes
  if (intpt.eq.1 .and. kfirst.eq.1)
c
  x call erhandler('usercr', 5000, 2,
  x 'ANSYS, INC.-supplied version of coding for USERCR
  x has been used.'
  x , 0.0d0, ' ')
c
  if (nuval*nintp.gt.840)
  x call erhandler('usercr', 5010, 4,
  x 'Maximum storage allowed by USVR has been exceeded.'
  x , 0.0d0, ' ')
c
c --- initialize creep strain increment in case on creep
  delcr = 0.0d0
c --- no creep if time is not moving
  if (timinc .le. 0.0d0 .and. timval .le. 0.0d0) fo to 999
c --- no creep if temperature is not defined
  temabs = tem + tofst
  if (temabs .le. 0.0d0) then
c
  call erhandler('usercr', 5020, 3,
  x 'Temperature= %G must be greater than zero for creep.'
  x , temabs, ' _
c
  go to 999
  endif

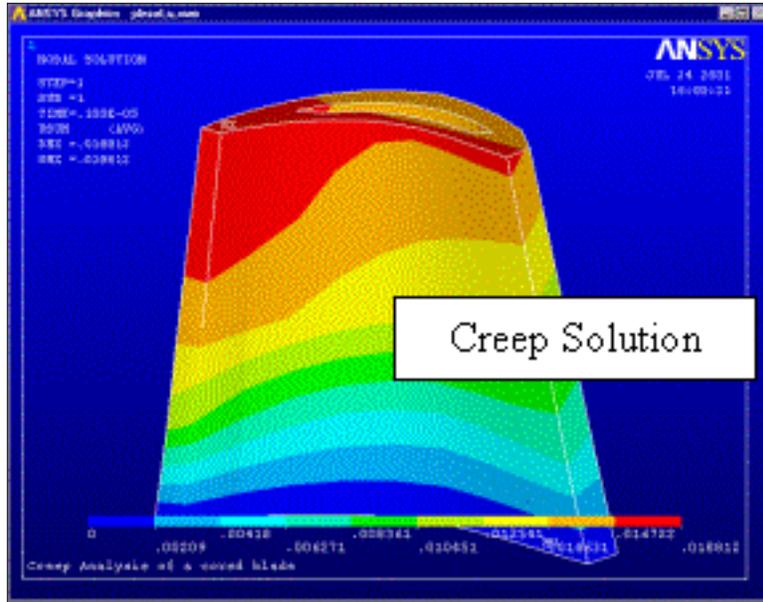
```

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<snip>



<snip>

```

c
c ***** define the equivalent strain and stress *****
c --- define the equivalent strain using the function egen
      epet = egen (nocomp,epel(1),posn)
c --- no cree if the strain is zero
      if (epet .eq. 0.0d0) go to 999
c --- define the stress
      sigen = e*epet
c
c ***** define the creep strain rate *****
c ***** normal beginning of user changes
c
c -PADT--- Put table values in constants for clarity
c
      c1 = proptb(1)
      alpha = proptb(2)
      dbin = proptb(3)
      vrbflg = proptb(5)
c
c -PADT--- If table item 4 (Q/k) is 0, set default of 8110
      if (proptb(4) .eq. 0) then

```



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```

    qovrk = 8110
  else
    qovrk = proptb(4)
  endif
<snip>

<snip>
c
c --- define the creep strain rate

    crprt = c1*(sinh(alpha*sigen)**dbltn)*dexp(-1*qovrk/temabs)

c ***** end of sample creep law
c ***** normal end of user changes
c
c ***** compute creep strain increments for each component *****
c
c --- define the creep strain increment from the rate
    delcr = crprt*timinc
c
c --- use prandtl-reuss relations to compute increments for each
    if (ncomp .eq. 1) then
      del(1) = delcr*epel(1)/epet
    else
      con = delcr/epet
      con = con/(2.0d0*(1.0d0+posn))
      call vmult (epel(4),del(4),ncomp-3,3.0d0*con)
      del(1) = con*(2.0d0*epel(1) - epel(2) - epel(3))
      del(2) = con*(2.0d0*epel(2) - epel(3) - epel(1))
      del(3) = con*(2.0d0*epel(3) - epel(1) - epel(2))
    endif
c
c --- update the strains
    call vapb1 (epcrp(1),del(1),ncomp)
    call vamb1 (epel(1),del(1),ncomp)
<snip>

<snip>
c -PADT--- Do Verbose I/O, if vrbflg eq 1
    if (intpt.eq.1 .and. vrbflg .eq. 1) then

```



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```

iott=6
write (iott,*) 'element number = ',elem,' gauss point 1'
write (iott,1) 'time=',timval,' stress=',sigen,' temp=',tem
write (iott,2) 'delcr=',delcr,' del=',del(1)
endif
c -PADT--- Do Summary I/O, if vrbflr eq 2
if (intpt.eq.1 .and. vrbflg .eq. 2) then
iott = 34
if (elem .eq. 1) then
epet = egen (ncomp,epel(1),posn)
write (iott,3) elem,timval,sigen,temabs,crprt,delcr,epet
endif
endif
c
999 return
1  format(2x,3(a,d15.8))
2  format(2x,2(a,d15.8))
3  format(2x,i6,6(',','g16.9))
end

```

## ANSYS Input File for USERCR

```

!ANSYS Input File for USERCR
! The ANSYS TB values are as follows:
! 1 C1
! 2 alpha-sub-1
! 3 n
! 4 Q/k (defaults to 8110)
! 5 Verbose output flag (1 = verbose output, 0 = standard output)
!
c1 = 9.62e4
a1 = 6e-4
n1 = 3.3
qovrk = 8110
vrbs = 2
TB,CREEP, 1
tbdata,6,100 !use user creep
tbdata,1,c1
tbdata,2,a1

```



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tbdata,3,n1  
tbdata,4,qovrk  
tbdata,5,vrbs

## UPFs: Lessons to Consider

- Always try APDL macros first
- Consider external commands to avoid custom versions of ANSYS
- You must have the proper FORTRAN compiler
  - See installation guide
- KISS
- Start by replicating the existing ANSYS functionality
  - Add your changes slowly
- Write and debug on simple models
- Don't be afraid of write(\*,\*)

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## FEA for .EDU

by [Rod Scholl](#), PADT

Finite Element Analysis (FEA) using a commercial code is becoming a more common skill for undergraduate mechanical engineers. Perhaps PADT can help you at your .edu institution by providing free course materials — and maybe more! The course material exchange program pools the resources of professors nationwide into an evolving curriculum that you can take advantage of. This and other free benefits may be available to you.

Please see the recent ANSYS [press release](#) for more information.



“They get younger every year!”

Drop us an [email](#) and we can discuss how we might help!



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## Upcoming Training at PADT

A complete schedule of the training courses for the next three months is shown below. [Learn more](#) about how the **Training Services** offered by PADT can save you time and money. Or, feel free to drop an e-mail to our training coordinator, [Ted Harris!](#)

Month	Start	End	Course	Description	Location
Apr '04	4/5	4/7	<a href="#">101</a>	<a href="#">Introduction to ANSYS, Part I</a>	Irvine, CA
	4/19	4/21	<a href="#">101</a>	<a href="#">Introduction to ANSYS, Part I</a>	Albuquerque, NM
	4/21	4/23	<a href="#">201</a>	<a href="#">Basic Structural Nonlinearities</a>	Tempe, AZ
May '04	5/3	5/5	<a href="#">101</a>	<a href="#">Introduction to ANSYS, Part I</a>	Tempe, AZ
	5/6	5/7	<a href="#">102</a>	<a href="#">Introduction to ANSYS, Part II</a>	Tempe, AZ
	5/10	5/11	<a href="#">100</a>	<a href="#">Engineering with Finite Element Analysis</a>	Tempe, AZ
	5/13	5/14	<a href="#">803</a>	<a href="#">Tcl/Tk for ANSYS</a>	Tempe, AZ
	5/19	5/20	<a href="#">203</a>	<a href="#">Dynamics</a>	Tempe, AZ
Jun '04	6/7	6/9	<a href="#">101</a>	<a href="#">Introduction to ANSYS, Part I</a>	Irvine, CA
	6/21	6/23	<a href="#">201</a>	<a href="#">Basic Structural Nonlinearities</a>	Tempe, AZ
	6/24	6/25	<a href="#">202</a>	<a href="#">Advanced Structural Nonlinearities</a>	Tempe, AZ
	6/28	6/29	<a href="#">301</a>	<a href="#">Heat Transfer</a>	Irvine, CA

Whether it s one of our regularly-scheduled classes at our training facility, or a customized class tailored to your specifications and location, you can be assured that the training you receive will have immediate, positive results on your capabilities in design, analysis, and product and process improvement.





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