

LS-DYNA® Database Binary Output Files

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INTRODUCTION

Three databases are discussed, these are:

- 1. State Database (G=ptf, default name d3plot and d3part)
- 2. Time History Database (F=thf, default name d3thdt)
- 3. Interface Force Database (S=iff, no default name, typically: intfor)

The purpose of this information is to give guidance on how to access and read the various databases.

The databases are written as word addressable fixed length binary files. The actual length depends on the amount of data saved, but will always be a multiple of 512 words (4 or 8 bytes each). Since it is likely that the database cannot be contained in a single file of length, FEMLEN, the data will spread over several files known as a family of files. Having a set of files enables them to be handled more easily than a single very large file. The root name for a family is the name of the first file member. Successive member names are compiled by appending a two or three digit number to the root name starting with 01, 02, through to 99, then 100 and ending with 999. Root names are limited to 75 characters. The original reason for a family of files was because the hard disks used for dyna3d runs could not cope with a single contiguous files large enough to contain all the data output. Subsequently, it has been found that splitting the output into separate files allows some unwanted data to be discarded and aids the copying, saving and movement of the data. Very large files can be impossible to transfer reliably over networks. Total output data can amount to several gigabytes or more depending on the model size. For Is-dyna runs with mesh adaptivity, the root name has a two letter appendage for each adapted mesh. Starting from 'aa' through 'az', then 'ba' through 'bz' and continuing up to 'zz', this gives a maximum of 676 possible adaptions. For example if the root name is 'd3plot' the subsequent files related to the original mesh are 'd3plot01', d3plot02, ..., after adaption the new mesh and undeformed geometry is put in 'd3plotaa' and the subsequent files for the new mesh are: 'd3plotaa01', d3plotaa02, ... The next adapted mesh will be in file 'd3plotab' and so on. A set of files at a particular adaption can be read separately by giving the root name with appendage, as the base file name.

For example command: 'lsprepost d3plot' will read in all the file with root name 'd3plot' While: 'lsprepost d3plotab' will read in only files have 'd3plotab' in the name.

LSPREPOST will read the binary databases separately or combined.

Eg: lsprepost d3plot, lsprepost d3thdt, lsprepost iffname, lsprepost d3plot h=d3thdt f=iffname

The file length used is set in the ls-dyna run as the default size of 7x512x512 words. The size can be changed on the command line with the 'x=factor' parameter giving a size of: factorx512x512 words.

If the initial data or state data is larger than the given file length, the data will automatically split across files. This condition is not desirable because it is not clear whether any non root file can be discarded without destroying the continuity of the data. Ls-dyna checks before writing to a file, to ensure that there is room left in the file to contain the data at a particular state time. If not, it closes the current family member and starts writing the state data in the next file member. The files are written with a block size of 512 words, and if the data does not complete the last block it is padded out. This means that files cannot be concatenated and read together. The word size is 4 bytes for the single precision version of ls-dyna and 8 bytes for the double precision version, unless 32bit ieee format is defined, see *DATABASE FORMAT, IBINARY.

FILE GENERAL STRUCTURE

The root file starts with a control words section, followed by node coordinates, then element connectivity for solids, thick shells, beams, and shells. Next are lists to reference the sequential internal numbering to the users number. State data is output next always starting with the time word. Data is of fixed length through the file members except where the mesh is adapted. The length of each area can be calculated from the information in the control words. The first file at adaption is like the root file in structure, so the new control words are used to recalculate the size of the subsequent data. The root file contains the initial data and also state data if there is room to write it. Further state data is written to the family members and each file will start with the time word provided data from the previous state did not overflow onto the file.

If the disk address, DA, of the data being written exceeds the maximum file length, then data is written into file number int(DA/FAMLEN) at location DA-FAMLEN*int(DA/FAMLEN). If the state length is greater than the remaining length, the disk address is increased to start the writing at the beginning of the next file.

STATE DATABASE (d3plot)

There are three sections in this database. The first contains 64 words of control information plus extensions. The second contains geometric information including the nodal coordinates and element connectivities and user numbering lists. The third section contains the results of the analysis at sequential output intervals. The output at a given time, called a state, contains a time word, global variables such as total energies and momenta for the whole model and each material (part), node data consisting of displacements, velocities, accelerations, and optionally temperatures, and finally element data that can include stresses and strains at integration points, and element deletion flags. The control data provides information about what is in the file and is used to calculate the various data length.

There are two other state database files, namely: d3drfl and d3part, these are similar to d3plot but contain less data. The dynamic relaxation file, d3drfl, provides the state at the end of the DR process, while d3part is state output for a reduced number of parts in the model.

CONTROL DATA

VALUE	#WORDS	ADDRESS	DESCRIPTION
Title	10	0	Model identification
Run time	1	10	time in seconds since 00:00:00 UTC, January 1, 1970
INUM (File ty)	pe) 1	11	d3plot=1
			1=d3plot, 2=d3drlf, 3=d3thdt, 4=intfor, 5=d3part

6=blstfor, 7=d3cpm, 8=d3ale, 11=d3eigv, 12=d3mode, 13=d3iter, 21=d3ssd, 22=d3spcm, 23=d3psd, 24=d3rms, 25=d3ftg, 26=d3acs

If > 1000, File type=INUM-1000

If INUM < 0 for d3plot and d3eigv then the data has be defined by the

*DATABASE_EXTENT_BINARY_COMP option.

DISK

all external(users) numbers (Node, Element, Part, and Rigid Surface Nodes) will be written in I8 format.

Length of arbitrary numbering array = NARBS * 8 bytes for single precision files.

Source version	1	12	ls-dyna version *1000000 + svn number
Release number	1	13	Release number in character*4 form
			50 for R5.0
			511c for R5.1.1c

Version	1	14	Code version, floating number, eg 960.0 it is used to distinguish the floating point format, like cray, ieee, and dpiece
NDIM	1	15	Number of dimensions (2 or 3). If 5 or 7 then an array of material types is read (MATTYP=1 is set in lsprepost), element connectivities are unpacked and NDIM=3. If 4 then element connectivies are unpacked in the DYNA3D database and NDIM is reset to 3. If >5 and < 8 then state data contains movement of rigid road surface. If NDIM=8 or 9, coordinates, velocities and accelerations are not output for rigid bodies nodes. At the end of each state motion data is output for each rigid body. See below for more details. If =9 rigid road surface data is included.
NUMNP	1	16	Number of nodal points
ICODE	1	17	Flag to identify finite element code=2: old DYNA3D,
			code=6: NIKE3D, LS-DYNA/3D, LS-NIKE3D
			database
NGLBV	1	18	Number of global variable to be read with each state
			NUMRW=number of rigid walls.
			NUMRBS=number of rigid body sets. = 6 + 7 * (NUMMAT8 + NUMMAT2 + NUMMAT4 + NUMATT+NUMRBS) + NUMRW * N N = 1 for DYNA3D and LS-DYNA3D N = 4 for LS-DYNA >= version 971
IT	1	19	Flag for temperatures = 0, none, = 1, read in a temperature for each node = 2, read temperature for each node and heat flux for each node. = 3, read thermal shell middle temperature, thermal shell inner temperature, thermal shell outer temperature, and heat flux for each node. Solid node temperatures are repeated +=10, read mass scaling value for each node
IU	1	20	Flag for current geometry (=1 or 0)
IV	1	21	Flag for velocities (=1 or 0). If <0 d3eigv file contains
			unscaled data.

IA NEL 8	1	22	Flag for accelerations (=1 or 0)
NEL8	1	23	Number of 8 node solid elements
			If NEL8 < 0, 2 extra nodes are output for ten node
			solids. Array is 2 * abs(NEL8), and is read after the
			arbitrary numbering arrays. NEL8 is also the total number of solid elements.
NUMMAT8	1	24	
NUMINIATO	1	2 4	Number of materials (parts) used by the 8 node solids
NUMDS	1	25	If < 0, Shell element data is output for 4 in-plane
			Gauss points. MAXINT will be 4 times number of
			points through the plane.
NUMST	1	26	=0 for d3plot/d3part
NV3D	1	27	Number of values in database for each solid element. =6*IOSOL(1)+IOSOL(2)+NEIPH If NV3D >= 8 * (6*IOSOL(1)+IOSOL(2)+NEIPH), each solid element has values at each Gauss point.
NEL2	1	28	Number of 2 node one-dimensional elements
NUMMAT2	1	29	Number of materials (parts) used by the 2 node 1D
			elements
NV1D	1	30	Number of values in database for each 1D element $= 6 + 5*BEAMIP + NEIPB*(3+BEAMIP)$.
NEL4	1	31	Number of four node shells (2D or 3D) elements.
NUMMAT4	1	32	Number of materials (parts) used by the 4 node 2D
			elements
NV2D	1	33	Number of values in database for each 2D element Are: <i>MAXINT</i> *(6*IOSHL(1)+IOSHL(2)+NEIPS)+8 *IOSHL(3)+4*IOSHL(4)+12*ISTRN
NEIPH	1	34	Number of additional values per integration point in a solid element, see NV3D. NEIPH is the number of values for history variables + total strains + plastic strains + thermal strains. ISTRN and IDTDT specifies which strains are present.
NEIPS	1	35	Number of additional values per integration point to be written into the type 6 database for shell elements.
MAXINT	1	36	Number of integration points dumped for each shell and the MDLOPT flag. The magnitude of <i>MAXINT</i> will be greater than or equal to 3. if MAXINT>=0, then MDLOPT=0 and <i>MAXINT</i> =MAXINT if MAXINT

			 MAXINT=abs(MAXINT) if MAXINT< -10,000, then MDLOPT=2 and MAXINT=abs(MAXINT)-10,000 MDLOPT controls the element deletion table (see below). This data allows deletion by nodes or elements.
EDLOPT	1	37	Element deletion flag (not standard) =xxx1 Solids deleted =xx1x Beams deleted =x1xx Shells deleted =1xxx Thick Shells deleted (Not used in LS-DYNA)
NMSPH	1	37	Number of SPH Nodes
NGPSPH	1	38	Number of SPH materials
NARBS	1	39	Additional storage required for arbitrary node and element numbering in type 6 database =0, Sequential numbering.
NELT	1	40	Number of 8 node thick shell elements. MAXINT*(6*IOSHL(1)+IOSHL(2)+NEIPS)+ 12*ISTRN
NUMMATT	1	41	Number of materials (parts) used for the 8 node thick shell element.
NV3DT	1	42	Number of values in database for each thick shell
IOSHL(1)	1	43	6 stress components flag. if 1000: IOSHL(1)=1, IOSOL(1)=1 if 999: IOSHL(1)=0, IOSOL(1)=1 else: IOSHL(1)=0, IOSOL(1)=0
IOSHL(2)	1	44	Plastic strain flag. if 1000: IOSHL(2)=1, IOSOL(2)=1 if 999: IOSHL(2)=0, IOSOL(2)=1 else: IOSHL(2)=0, IOSOL(2)=0
IOSHL(3)	1	45	Shell force resultants flag, if 1000 =1 else =0
IOSHL(4)	1	46	Shell thickness, energy+2 others, if 1000 =1 else =0
IALEMAT	1	47	Size of array containing solid element parts numbers used as ALE material
NCFDV1	1	48	Bit flags for CFD nodal values, or if = 67108864, then state contains Multi-Solver extra data – see formats descriptions below.

NCFDV2	1	49	Further bit flags for CFD nodal values or if MS extra data, then value equals number of data domains.	
NADAPT	1	50	Number of adapted element to parent pairs (not implemented)	
NMMAT	1	51	Total number of materia DYNA/3D	als (parts) – not set in LS-
NUMFLUID	1	52	volume fractions output for the dominant group. species mass for each g vf1, vfn, dvf flag, m after the location for pla history variables are wr	uid groups. Fluid density and as history variables, and a flag. If negative multi-material roup is also output. Order is: rho, 1, mn. Density is at position 8 astic strain. Any element material itten before the Ale variables, ins components after these if
INN	1	53	Invariant node numbering See INN in card *CON*	ng fore shell and solid elements ΓROL_ACCURACY
NPEFG	1	54		hod data sets. When the seventh Discrete Element Sphere (DES) NPEFG/1000000 = 1.
NEL48	1	55	extra nodes are output f * NEL48 long, and is re	Is. Internal element id and 4 for each 8 node shell. Array is 5 ead before the Header, Part and er the extra nodes for 10 nodes
IDTDT	1	56	•	n the database, these are: unit, and so on. Extract the digit from following types:
			IDTDT / 1 = 1:	An array of dT/dt values of length NUMNP. Array is written after node temperature arrays. (xxxx1)
			IDTDT! 10 = 1:	An array of residual forces of length 3*NUMNP followed by residual moments of length 3*NUMNP. This data is written after node temperatures or dT/dt values if there are output. (xxx1x)

IDTDT! 100 = 1: Plastic strain tensor is written

for each solid and shell after standard element data. For solids (6 values) and for shells (6 x 3 = 18 values), at the lower, middle and upper integration location. (xx1xx)

IDTDT! 1000 = 1: Thermal strain tensor is written

after standard element data. For solid (6 values) and shell (6 values) and after the plastic strain tensor if output. (x1xxx)

IDTDT ! 10000 = 1: If database contains thermal or

plastic strains, IDTDT>100, then this is the value of ISTRN,

= 0 or 1. (1xxxx)

EXTRA 1 57 Additional number of control words. If > 0, there are
EXTRA control words after the first 64 words.

WORDS 6 58-63 Used by D3THDT and INTFOR

The value of ISTRN must be computed if IDTDT<100, because then it is not output in the control data. For this case only, this is the rule for computing it.

ISTRN can only be computed as follows and if NV2D > 0.

If NV2D-MAXINT*(6*IOSHL(1)+IOSHL(2)+NEIPS)+8*IOSHL(3)+4*IOSHL(4) > 1

Then ISTRN = 1, else ISTRN = 0

If ISTRN=1, and NEIPH>=6, last the 6 additional values are the six strain components.

Or NELT > 0

If NV3DT-MAXINT*(6*IOSHL(1)+IOSHL(2)+NEIPS) > 1

Then ISTRN = 1, else ISTRN = 0

If (EXTRA > 0) Extra control words are written after the first 64 words (This is currently set to 64 and 12 are used from September 2014, remaining are set to zero)

DISK

NEL20	1	64	Number of 20 node Solid Hexahedron Elements. Internal element id and 12 extra nodes are output for 20 node solids. Array is 13 * NEL20 long, and is read before Header, Part and Contact details, and after the extra nodes for 8 node shells, if they exist.
NT3D	1	65	Number of Thermal Element Variables, Data is NT3D * NEL8 and is output after the thermal nodal data.
NEL27	1	66	Number of 27 node Solid Hexahedron Elements.
NEIPB	1	67	Number of requested material history variables per integration point for beam elements.
NEL21P	1	68	Number of 21 node Solid Pentahedron Elements
NEL15T	1	69	Number of 15 node Solid Tetrahedron Elements.
SOLENG	1	70	If SOLENG>0, this is the position where to find
			internal energy density for solid elements in solid
			element integration point data.
NEL20T	1	71	Number of 20 node Solid Tetrahedron Elements.
NEL40P	1	72	Number of 40 node Solid Pentahedron Elements.
NEL64	1	73	Number of 64 node Solid Hexahedron Elements.
QUADR	1	74	Flag for output of state data for Quadratic elements
			=1 for full element connectivity, =2 for full
			connectivity and state data at each integration point.
CUBIC	1	75	Flag for output state data for Cubic elements
TSHENG	1	76	If TSHENG > 0 , this is the position where to find
			internal energy density for thick shell elements.
NBRANCH	1	77	Number of post branches
PENOUT	1	78	0/1/2. Flag for contact penetration output. =1: output
			max contact penetration, x, y, z for each node
			(3*NUMNP values). =2, in addition to absolute
			penetrations, also output relative (to max possible) x,
			y, z penetrations. (3*NUMNP values). Data written
			after residual forces and moments.
ENGOUT	1	79	0/1. Flag for output contact energy density as a node
			scalar field. Data written after contact penetrations.

These flags are set in *DATABASE_EXTENT_BINARY card 4 columns 4 and 5.

HIGHER SOLID ELEMENT PART DATA

If QUADR or CUBIC is > 0

part data is output for each part defined with higher order elements

NPART	1	63+EXTRA	Number of higher order element parts
repeat for each parts	s:		
PID	1	64+EXTRA	Part internal id.
MID	1	65+EXTRA	Part material id.
EOS	1	66+EXTRA	Material equation of state.
EFORM	1	67+EXTRA	Element formation id.
NMNP	1	68+EXTRA	Number of nodes in element.
NGP	1	69+EXTRA	Number of integration points in element.
LENGP	1	70+EXTRA	Number of variables per integration point.
NHISV	1	71+EXTRA	Number of history variables.
ISTRN	1	72+EXTRA	Strain flag indicates output of strain tensor.

MATERIAL TYPE DATA

The material section contains the material type numbers. This section is skipped if MATTYP is zero.

This data is required because those shell elements that are in a rigid body have no element data output in the state data section. The normal length of the shell element state data is:

NEL4 * NV2D, when the MATTYP flag is set the length is: (NEL4 – NUMRBE) * NV2D.

When reading the shell element data, the material number must be check against IRBRTYP list to find the element's material type. If the type = 20, then all the values for the element to zero.

This option is set in *DATABASE_EXTENT_BINARY, with DCOMP=2

VALUE	LENGTH	DESCRIPTION
NUMRBE	1	Number of rigid body shell elements.
NUMMAT	1	Number of materials in the database.
IRBTYP	NUMMAT	Material type numbers

FLUID MATERIAL ID DATA

The fluid material section contains the material numbers for solid elements that are used to define an Euler grid or Arbitrary Lagrangian Euler mesh. This section is skipped if IALEMAT is zero.

VALUE LENGTH DESCRIPTION

FLUIDID IALEMAT Fluid material number used in solid element mesh

SMOOTH PARTICLE HYDRODYNAMICS ELEMENT DATA FLAGS

This section is only output if NMSPH > 0. The section is a list of flags to indicate what SPH data is output for each SPH node/element. The first number is the length in words for this array, currently = 11.

SPH elements are centered at nodes, and cover a spherical volume defined by the radius of influence. They do not have a connection with other SPH elements. They should be displayed as a dot or a spherical surface, with radius scaling to reduce the size and enable each element to be distinguishable.

As follows:

```
isphfg(1) = 11 - length of sph flags array
isphfg(2) = 1 - radius of influence
isphfg(3) = 1 - pressure in particle
isphfg(4) = 6 - 6 true stress components
isphfg(5) = 1 - plastic strain, > 0.0 if effective stress exceeds yield strength
isphfg(6) = 1 - density of particle material
isphfg(7) = 1 - internal energy (strain)
isphfg(8) = 1 - number of neighbors affecting particle
isphfg(9) = 12 - 6 strain (ex, ey, ez, exy, eyz, exz) and 6 strain rate (erx, ery, eryz, eryz, eryz, eryz)

if negative, true strain components
isphfg(10)=1 - mass of element (>= ls971)
isphfg(11)=1 - max number of sph history variables.
```

If any value of isphfg(2) through isphfg(11) = 0, then the particular data item is not output for the particle. To calculated the size of data add the isphfg values from isphfg(2) through isphfg(11) plus one. One value is always output which is the material number as a floating point number for each particle.

If this value is negative then the particle has been deleted from the model.

Note: it is possible a SPH element could be deleted, or be non active in the initial states, and become active in later states.

Full output for each particle is:

```
mat#, radius, pressure, {sx, sy, sz, sxy, syz, sxz} ps, rho, ie, nn, {ex, ey, ez, exy, eyz, exz, erx, ery, erz, erxy, eryx, erxz}, mass, hv1 ... hvn.
```

```
NUM\_SPH\_VARS = 1 + sum of isphfg(i), i=2 to isphfg(1)
```

Hence, total size is 20 + the total number of history variables.

When a particle is deleted from the model, data is still output for it because the length of data must always be the same for each state.

PARTICLE DATA (NPEFG > 0)

DES Control Block – see description below

Control Block

If NPEFG > 0 and NPEFG < 10000000, airbag particles are output

The first three digits of NPEFG are the number of airbags in the database = NPARTGAS

NPARTGAS = NPRFG % 1000

SUBVER = NPEFG / 1000

In the extended control block:

The first four words in the block are:

1. NGEOM number of geometry variables

2. NVAR number of state variables

3. NPART number of particles

4. NSTGEOM number of state geometry variables

If SUBVER == 4

5. NCHAMBER number of chambers

NLIST = NGEOM + NVAR + NSTGEOM

NLIST words of output for variables listed to define the type of each variable, =1 for integer and 2= for floating point

8 * NLIST words of 8 character variable names (each integer word is an ascii character).

GEOMETRY DATA

The geometry section contains the nodal coordinates and the element connectivities. The ordering of the nodal points is the same as the ordering of the nodal data in the state data that follows. If NDIM=3 the connectivities are assumed to be packed with 3 integers per word, if NDIM>3, then connectivities are <u>not pack</u>, (the default for LS-DYNA, LS-DYNA3D and LS-NIKE3D. The order of the elements are 3, 2, and 1 dimensional elements if the database is ICODE=2 or 6.

VALUE	LENGTH	DESCRIPTION
X(3,1)	NDIM*NUMNP	Array of nodal coordinates X1,Y1,Z1, X2,Y2,Z2, X3,Y3,Z3,,Xn,Yn,Zn
IX8(9,1)	9*NEL8	Connectivity and material number for each 8 node solid element. Also, corner nodes for other > 8 node solid elements. Pentahedron is: 1,2,3,4,5,5,6,6, and Tetrahedron is: 1,2,3,4,4,4,4,4
If $NEL8 < 0$	2*abs(NEL8)	Extra nodes for ten node solids – not read here.
IXT(9,1)	9*NELT	Connectivity and material number for each 8 node thick shell element.
IX2(6,1)	6*NEL2	Connectivity, orientation node, two null entries, and the material number for each 2 node beam element. For some beam types the last two number contain the beam type and length to width ratio * 100 and length to height ratio * 100 type = $ix2(5,*)$ & $0x3F$ width = $0.01 * length / (ix2(5,*)>>6$ height = $0.01 * length / ix2(6,*)$ Third node (orientation) may be > 1e9 Contain flag 1e9 to indicate a spot weld.
IX4(5,1)	5*NEL4	Connectivity and material number for each 4 node shell element

Note the node numbers are the LS-DYNA internal numbers for nodes, these will be the same as the user's numbers if NARBS = 0, otherwise, the arbitrary number lists are used to find the user's numbers, similarly, for element numbers and material numbers.

USER MATERIAL, NODE, AND ELEMENT IDENTIFICATION NUMBERS

Skip this section if NARBS (disk address 39) is zero. The user node and element numbers must be in ascending order. *It assumed that if this option is used all node and element data anywhere in the databases is in ascending order in relation to the user numbering*. Read in NARBS words and decipher as indicated below.

For sequential material/part numbering, the total length of data is:

NARBS=10+NUMNP+NEL8+NEL2+NEL4+NELT+

3*NMMAT: these numbers are not used

Only the first 10 control words are read in and used. The other 6 words are only output by ls-dyna when NSORT < 0.

For arbitrary material numbering (NSORT < 0)

NARBS=16+NUMNP+NEL8+NEL2+NEL4+NELT+3*NMMAT

Where material numbers are not in ascending order.

For this case all 16 control words are read in.

VALUE	LENGTH	DESCRIPTION	
NSORT	1	Pointer to arbitrary node numbers in LS-DYNA code. If < 0 , it flags that arbitrary material identification numbers are also used.	
NSRH	1	Pointer to arbitrary solid element numbers in LS-DYNA code: =NSORT+NUMNP	
NSRB	1	Pointer to arbitrary beam element numbers in LS-DYNA code: =NSRH+NEL8	
NSRS	1	Pointer to arbitrary shell element numbers in LS-DYNA code: =NSRB+NEL2	
NSRT	1	Pointer to arbitrary thick shell element numbers in	
		LS-DYNA code: =NSRS+NEL4	
NSORTD	1	Number of nodal points	
NSRHD	1	Number of 8 node solid elements	
NSRBD	1	Number of 2 node beam elements	
NSRSD	1	Number of 4 node shell elements	
NSRTD	1	Number of 8 node thick shell elements	

1	Pointer to an array in the LS-DYNA code that list the material ID's in ascending order.	
1	Pointer to an array in the LS-DYNA code that gives the material ID's in the actual order that they are defined in the user input.	
1	Pointer to an array in the LS-DYNA code that gives the location of a member in the array originating at NSRMU for each member in the array starting at NSRMA.	
1	Total number of materials (parts) Total number of nodal rigid body constraint sets	
1	Total number of materials	
1		
NSORTD	Array of user defined node numbers	
NSORTH	Array of user defined solid element numbers	
NSORTB	Array of user defined beam element numbers	
NSORTS	Array of user defined shell element numbers	
NSORTT	Array of user defined thick shell numbers	
NMMAT	Ordered array of user defined material (part) ID's	
NMMAT	Unordered array of user material (part) ID's	
NMMAT	Cross reference array	
	1 1 1 NSORTD NSORTH NSORTB NSORTS NSORTT NMMAT NMMAT	

RIGID BODY DESCRIPTION for NDIM=8,9

NRIGID 1		Number of rigid bodies			
For each rigid body:					
MRIGID	1	Rigid body part internal number			
NUMNODR	1	Number of nodes in rigid body			
NODER	NUMNODR	Internal node number of rigid body			
NUMNODA	1	Number of active (not rigid) nodes			
NODEA	NUMNODA	Internal node numbers of active nodes.			

Active node coordinates, velocities and accelerations are output as usual. Rigid body motion data is described in the state section below.

ADAPTED ELEMENT PARENT LIST (not implemented)

List of element id pairs for H-type shell element adaptivity.

Length of data is 2 * NADAPT, pairs are element number and element parent number

SMOOTH PARTICLE HYDRODYNAMICS NODE AND MATERIAL LIST

If NMSPH > 0 List of sph node and its material number

Length of data 2 * NUMSPH

PARTICLE GEOMETRY DATA (NPEFG > 0)

DES Control words (NPEFG/10000000 == 1) – see description below

NPARTGAS blocks of NGEOM data to describe the geometry for each airbag:

- 1. first particle ID for the airbag
- 2. number of particles in the airbag
- 3. ID for the airbag
- 4. number of gas mixtures in the airbag

If NGEOM == 5

5. number of chambers

RIGID ROAD SURFACE DATA

If NDIM > 5

NNODE Number of nodes in road surface

NSEG Total number of 4 noded road surface segments

NSURF Number of road surfaces

MOTION Flag to indicate motion data is output for each state

NODEID NNODE list of IDs

SURFNODE XYZ Coordinate for each node

Lists of 4 node segments for each surface

SURFID Surface ID Number

SURFNSEG Number of segments in surface

SURFSEGS SURFNSEG of 4 node ids for each segment

Length of data = 4 + NNODE + 3 * NNODE + NSURF * (2 + 4 * SURFNSEG)

EXTRA 2 NODE CONNECTIVITY FOR 10 NODE TETRAHEDRON ELEMENTS

(only if NEL8 < 0)List of extra nodes for each 10 node tetrahedron element, 2 * abs(NEL8). Any 8 node solids have these two nodes set to zero.

EXTRA 4 NODE CONNECTIVITY ARRAY FOR 8 NODE SHELL ELEMENTS

(only if NEL48 > 0)

NEL48 number of 8 node shells

List of extra nodes for each 8 node shell element, 5 * NEL48:- element internal number and 4 extra nodes.

EXTRA 12 NODE CONNECTIVITY ARRAY FOR 20 NODE HEXAHEDRON ELEMENTS

(only if EXTRA > 0 and NEL20 > 0)

NEL20 number of 20 node solids

List of extra midside nodes for each 20 node solid elements, 13 * NEL20:- element internal number and 12 extra nodes for each element. Corner nodes are taken from NEL8 list.

27 NODE CONNECTIVITY ARRAY FOR 27 NODE HEXAHEDRON ELEMENTS

(only if EXTRA > 0 and NEL27 > 0 and QUADR > 0)

NEL27 number of 27 node solids

List of 27 nodes for each 27 node solid elements, 28 * NEL27:- element internal number and 27 nodes for each element. Ls-dyna R10 outputs 19 nodes for each element to be combined with the 8 corner nodes already output.

21 NODE CONNECTIVITY ARRAY FOR 21 NODE PENTAHEDRON ELEMENTS

(only if EXTRA > 0 and NEL21P > 0 and QUADR > 0)

NEL21P number of 21 node solids

List of nodes for each 21 node solid element, 22 * NEL21P:- element internal number and 21 nodes for each element.

15 NODE CONNECTIVITY ARRAY FOR 15 NODE SOLID ELEMENTS

(only if EXTRA > 0 and NEL15T > 0 and QUADR > 0)

NEL15T number of 15 node solids

List of nodes for each 15 node solid elements, 8 * NEL15T:- element internal number and 15 nodes for each element.

20 NODE CONNECTIVITY ARRAY FOR 20 NODE TETRAHEDRON ELEMENTS

(only if EXTRA > 0 and NEL20T > 0 and CUBIC > 0)

NEL20T number of 20 node solids

List of nodes for each 20 node solid elements, 21 * NEL20T:- element internal number and 20 nodes for each element.

40 NODE CONNECTIVITY ARRAY FOR 40 NODE PENTAHEDRON ELEMENTS (only if EXTRA > 0 and NEL40P > 0 and CUBIC > 0)

NEL40P number of 40 node solids

List of nodes for each 40 node solid elements, 41 * NEL40P:- element internal number and 20 nodes for each element.

64 NODE CONNECTIVITY ARRAY FOR 64 NODE HEXAHEDRON ELEMENTS (only if EXTRA > 0 and NEL64 > 0 and CUBIC > 0)

NEL64 number of 64 node solids

List of nodes for each 64 node solid elements, 65 * NEL64:- element internal number and 64 nodes for each element.

HEADER, PART & CONTACT INTERFACE TITLES

At the end of the first binary files, eg d3plot, the part and model titles are appended. If the model input includes *DATABASE_BINARY_D3PROP, all the d3prop part data is included.

At the end of the first interface force file, titles and contact id are appended.

This extra data is written at the end of the following files: d3plot, d3part and intfor files, and the header and part titles are written directly after the EOF (= -999999.0) marker.

Header output

NTYPE 1 entity type = 90000

HEAD 18 Header title (72 characters)

For the interface force file (intfor), header and contact titles are written at the end of first file after the EOF (= -999999.0) marker

Part title output

Value	Length	Description		
NTYPE	1	entity type = 90001		
NUMPROP	1	number of parts		

For NUMPROP parts:

IDP 1 part id

PTITLE 18 Part title (72 characters)

For the interface force file (intfor), header and contact titles are written at the end of first file after the EOF (= -999999.0) marker.

Contact title output

NTYPE 1 entity type = 90002 NUMCON 1 number of contacts

For NUMCON contacts:

IDC 1 contact id

CTITLE 18 Contact title (72 characters)

Header output

NTYPE 1 entity type = 90000

HEAD 18 Header title (72 characters)

ICFD Part title output

Value Length Description

NTYPE 1 entity type = 90020 NUMPROP 1 number of parts

CESE: mechanics solid surface part title output

Value Length Description

NTYPE 1 entity type = 90021 NUMPROP 1 number of parts

For NUMPROP parts:

IDP 1 part id

PTITLE 18 Part title (72 characters)

The d3prop data is written to the d3plot file only if it is requested.

D3PROP output

Values Length Description

NTYPE 1 entity type = 900100 NLINE 1 number of keyword lines

For NLINE keyword lines:

KEYWORD 20 keyword line (80 characters)

DESCRIPTION OF BINARY FILE TYPES

Control word 11

File type:

1=d3plot plot file of model and state data

2=d3drlf plot file of model and state data from a dynamic relaxation analysis

3=d3thdt time history plot file for a set of nodes and elements

4=intfor plot file of contact interfaces

5=d3part plot file of model and state data for a set of parts

6=blstfor plot file for a blast wave analysis

7=d3cpm

8=d3ale plot file for ale fluid-structure interface

or fsifor

11=d3eigv plot file for an eigen value analysis

12=d3mode

13=d3iter

21=d3ssd plot file for steady state dynamic response.

22=d3spcm plot file for response spectrum analysis.

23=d3psd plot file for power spectral density of response, in random vibration.

24=d3rms	plot file for root mean square of response, in random vibration.
25=d3ftg	plot file for random fatigue analysis.

26 12	1 4 6.1 6	C	1 .	' PPA 1 '
26=d3acs	nior rue ro	or treamency	domain	acoustic FEM analysis
_ 0	prot mie	or moderno	COLLICIA	acoustic I Bill and Jois

EXTRA DATA TYPES (Output for Multi-Solver Analysis)

If NCFDV1 = 67108864, then NCFDV2 will be the number of additional datasets from different solver-mesh combinations that are included in the d3plot file. One of each of the solver-mesh combinations listed below can be among the NCFDV2 datasets. Currently defined solver-mesh combinations follow.

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of volume vars output: nvolvar_pfem
first volume variable ID: ID 1
...
last volume variable ID: ID nvolvar_pfem
number of PFEM parts
first internal part ID: partID 1
...
last internal part ID: partID nPFEM_parts
first user part ID: user_partID 1
...
last user part ID: user_partID nPFEM_parts
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output:

nsurfvar_pfem
first surface variable ID:
last surface variable I
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
first internal part ID: partID 1
...
last internal part ID: partID nCESE_parts
first user part ID: user_partID 1
...
last user part ID: user_partID nCESE parts
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output:

nsurfvar_cese
first surface variable ID:

last surface variable ID:

ID 1

...

last surface variable ID:

ID nsurfvar_cese

number of CESE surface parts
first internal part ID:

last internal part ID:

partID 1

...

partID nCESE_surfparts
first user part ID:

user_surfpartID 1

...

last user part ID:

user_surfpartID nCESE_surfparts
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of volume vars output:

nvolvar_cese
first volume variable ID:

last volume variable ID:

ID nvolvar_cese

number of CESE parts
first internal part ID:

last internal part ID:

partID 1

partID nCESE_parts
first user part ID:

user_partID 1

user_partID 1
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output:
nsurfvar_cese
first surface variable ID:
last surface variable ID:
ID nsurfvar_cese
number of CESE surface parts
first internal part ID:
partID 1
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of volume vars output:

nvolvar_cese
first volume variable ID:

last volume variable ID:

ID nvolvar_cese

number of CESE parts
first internal part ID:

last internal part ID:

partID 1

partID nCESE_parts
first user part ID:

user_partID 1

user_partID 1
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output:
first surface variable ID:
last surface parts
first internal part ID:
last internal part ID:
last user surfpartID nCESE surfparts
user_surfpartID nCESE surfparts
```

In this domain, the variables are defined at the element centroid.

```
solver and domain ID:

size of each variable component: numelh_cese
number of nodes:
number of elements:
numelh_cese
number of elements:
numelh_cese
user node numbers:
array of nodal coordinates:
element connectivity:
number of output vars:
nv_cese_cfd(9, numelh_cese)
number of output vars:
nv_cese_cfd_ele
first variable ID:
last variable ID:
number of CESE parts
first internal part ID:
partID 1
```

In this domain, the variables are defined at the surface element centroid.

```
solver and domain ID:

size of each variable component: numelsurf_cese
number of nodes:
number of elements:
user node numbers:
array of nodal coordinates:
element connectivity:
number of output vars:
number of output vars:
nv_cese_cfd(3, numnp_cese)
number of output vars:
nv_cese_cfd_ele
first variable ID:
lD nv_cese_cfd_ele
number of CESE surface parts
first internal part ID:
last internal part ID:
last user part ID:
user_surfpartID ncese_surfparts
for the first CESE surface element

user_surface element
for the last CESE surface element
```

In this domain, the variables are defined at the 2D element centroid.

```
solver and domain ID:

size of each variable component:
numel2d_cese
number of nodes:
number of elements:
user node numbers:
array of nodal coordinates:
element connectivity:
number of output vars:
first variable ID:
number of CESE parts
first internal part ID:
last internal part ID:
partID ncese_parts
first user part ID:
user_partID ncese_parts
user_partID ncese_parts
for the first 2D CESE element

cuser_element number:

size of each variable component:
numnel2d_cese
numnp_cese
numel2d_cese
numel2d_cese)
numel2d_cese)
numel2d_cese
```

In this domain, the variables are defined at the 2D surface element centroid.

solver and domain ID:
size of each variable component: numsurf2Dele_cese
number of nodes:
number of elements:
user node numbers:
array of nodal coordinates:
element connectivity:
number of output vars:
number of output vars:
nv_cese_cfd_ele
first variable ID:
last variable ID:
last variable ID:
last internal part ID:
last internal part ID:
last user part ID:
user_surfpartID ncese_surfparts
for the first CESE surface element
user_element number:

for the last CESE surface element

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In this domain, the variables are defined at the 2D axisymmetric element centroid.

solver and domain ID:

size of each variable component: numel2daxi_cese
number of nodes:
number of elements:
user node numbers:
array of nodal coordinates:
element connectivity:
number of output vars:
first variable ID:
number of CESE parts
first internal part ID:
last internal part ID:
partID ncese_parts
first user part ID:
user_partID ncese_parts
user element number:

for the last 2D CESE element

cese
numnp_cese
numnp_cese
numnp_cese)
numnp_cese)
numel2daxi_cese)
numel2daxi_cese)
numnp_cese
nump_cese

In this domain, the variables are defined at the 2D axisymmetric surface element centroid.

```
user node numbers:
                              nodes cese cfd(numnp cese)
array of nodal coordinates: x_cese_cfd(3, numnp_cese)
                           ix2_cese_cfd(3,
nv_cese_cfd_ele
element connectivity:
                              ix2 cese cfd(3, numsurfel2daxi cese)
number of output vars:
                              ID^{-}1
first variable ID:
last internal part ID:
                            partID ncese_surfparts
first user part ID:
                              user surfpartID 1
last user part ID: user_surfpartID ncese_surfparts user element number: for the first CESE surface element
user element number:
                      for the last CESE surface element
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output: nsurfvar_mech first surface variable ID: ID 1
...
last surface variable ID: ID nsurfvar_mech
number of MECH surface parts nMECH_surfparts first internal part ID: partID 1
...
last internal part ID: partID nMECH_surfparts first user part ID: user_surfpartID 1
...
last user part ID: user_surfpartID nMECH surfparts
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output: nsurfvar_mech
first surface variable ID: ID 1
...
last surface variable ID: ID nsurfvar_mech
number of MECH 2D surface parts first internal part ID: partID 1
...
last internal part ID: partID nMECH2D_surfparts
first user part ID: user_surfpartID 1
...
last user part ID: user_surfpartID 1
...
last user part ID: user surfpartID nMECH2D surfparts
```

In this domain, the variables are defined on structural solid elements.

```
solver and domain ID:

size of each variable component: nip_solid_em * numelh
number of nodes:
numnp
number of elements:
numelh
number of integration points:
numelh
number of output vars:
nv_em_solid_em
first variable ID:
ID 1

...
last variable ID:
ID nv_em_solid_integ
```

In this domain, the variables are defined on structural thick shell elements.

```
solver and domain ID:

size of each variable component: nip_tshell_em * numelt
number of nodes:
numnp
number of elements:
number of integration points:
number of output vars:
first variable ID:

last variable ID:

ID nv_em_tshell_integ
```

In this domain, the variables are defined on structural thin shell elements.

```
solver and domain ID:

size of each variable component: nip_shell_em * numels
number of nodes:
numnp
number of elements:
number of integration points:
number of output vars:
number of output vars:
first variable ID:

last variable ID:

ID nv_em_shell_integ

ID nv_em_shell_integ
```

In this domain, the variables are defined at the centroids of structural solid elements.

```
solver and domain ID: EM_FEMSTER_SOLID_CENTROID size of each variable component: numelh number of nodes: numnp number of elements: numelh number of output vars: nv_em_solid_cent first variable ID: ID 1 ... last variable ID: ID nv_em_solid_cent
```

In this domain, the variables are defined at the centroids of structural thick shell elements.

In this domain, the variables are defined at the centroids of structural thin shell elements.

In this domain, the variables are defined at the integration points of structural beam elements.

```
solver and domain ID: EM_FEMSTER_BEAM_INTEG_PTS size of each variable component: numelb number of nodes: numnp number of elements: numelb number of output vars: nv_em_beam_cent first variable ID: ID 1 ... last variable ID: ID nv_em_beam_cent
```

In this domain, the variables are defined at the mesh nodes.

```
solver and domain ID:

size of each variable component: nip_air_em * number of nodes:
number of elements:
number of integration points:
nuser node numbers:
array of nodal coordinates:
element connectivity:
number of output vars:
first variable ID:

EM_FEMSTER_AIR
nip_air_em * numelh_air_em
numnp_air_em
numnp_air_em
nip_air_em
nodes_air_em(numnp_air_em)
ix8_air_em(3, numnp_air_em)
ix8_air_em(8, numelh_air_em)
number of output vars:
first variable ID:

ID nv_em_air_integ
```

In this domain, the variables are defined at the nodes of the implied rectangular mesh.

```
solver and domain ID:
                                              RECT AIR EM NODE
size of each variable component: nx rect air em * ny rect air em
                                              * nz rect air em
number of x nodes:
                                             nx_rect_air_em
                                      nx_rect_air_em
ny_rect_air_em
nz_rect_air_em
xmin_rect_air_em
ymin_rect_air_em
zmin_rect_air_em
ymax_rect_air_em
ymax_rect_air_em
zmax_rect_air_em
zmax_rect_air_em
nv_em_air_nd
ID 1
number of y nodes:
number of z nodes:
minimum x coordinate:
minimum y coordinate:
minimum z coordinate:
maximum x coordinate:
maximum y coordinate:
maximum z coordinate:
number of output vars:
first variable ID:
                                            ID 1
```

```
last variable ID: ID nv em air nd
```

In this domain, the variables are defined on faces of structural elements.

```
solver and domain ID:
                                   EM FEMSTER BEM
size of each variable component: nip bem em * nfaces bem em
number of nodes: numnp_bem_em number of elements: nfaces_bem_em number of integration points: nip_bem_em
number of BEM parts:
                                 em numPartBem
flag for BEM mesh:
                                  nBEMflag
        (first bit =0 if no motion,=1 if motion)
        (second bit =0 if no edge domain, =1 if edge domain)
        (third bit =0 if no node domain, =1 if node domain)
user node numbers:
                                  nodes bem em (numnp bem em)
array of nodal coordinates: x_bem_em(3, numnp_bem_em)
element connectivity:
                                  ix4 bem em(5, nfaces bem em)
if (second bit(nBEMflag) = 1)
 endif
 f (third bit(nBEMflag) = 1)
node element connectivity number of node domains node domain array nodeDomain(numnp_bem_em)
nodex4_bem_em(5, nfaces_bem_em)
if (third bit(nBEMflag) = 1)
endif
number of output vars: nv_em_bem_integ
                                  ID^{-}1
first variable ID:
last variable ID:
                                   ID nv em bem integ
```

In this domain, the variables are defined at the nodes of the mechanics structural mesh.

```
solver and domain ID:

number of nodes:

numnp
size of each variable component:
numnp_em
node list (internal #s):
numnp_em
number of output vars:
first variable ID:

last variable ID:

ID nv node em
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

solver and domain ID: DUALCESE3D ADAPT last volume variable ID: ID nvolvar dualcese number of DUALCESE3D parts nDUALCESE3D_parts first internal part ID: partID 1 last internal part ID: partID nDUALCESE3D_parts first user part ID: user partID 1 last user element number: for the last DUALCESE3D ADAPT element

Notes:

(1) The user elements numbers only refer to the unrefined mesh elements. As no elementcentered data is output, these are only provided for user reference.

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

solver and domain ID: DUALCESE3D SURF ADAPT number of surface vars output: nsurfvar_dualcese ID 1 first surface variable ID: last surface variable ID: ID nsurfvar dualcese first internal part ID: partID 1 user surfpartID nDUALCESE3D surfsetIDs last user part ID:

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

solver and domain ID: DUALCESE2D ADAPT number of volume vars output: nvolvar_dualcese first volume variable ID: last internal part ID:

partID nDUALCESE parts

```
first user part ID:

...

last user part ID:

user_partID 1

user_partID nDUALCESE_parts

initial number of elements:

nel

first user element number:

for the first DUALCESE2D_ADAPT element

...

last user element number:

for the last DUALCESE2D ADAPT element
```

(1) The user elements numbers only refer to the unrefined mesh elements. As no element-centered data is output, these are only provided for user reference.

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output: nsurfvar_dualcese first surface variable ID: ID 1
...
last surface variable ID: ID nsurfvar_dualcese
number of DUALCESE surf. setIDs nDUALCESE_surfsetIDs first internal part ID: partID 1
...
last internal part ID: partID nDUALCESE_surfsetIDs first user part ID: user_surfpartID 1
...
last user part ID: user_surfpartID nDUALCESE_surfsetIDs
```

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of volume vars output:
first volume variable ID:
last volume var
```

Notes:

(1) The user elements numbers only refer to the unrefined mesh elements. As no element-centered data is output, these are only provided for user reference.

For the following domain, the mesh can be completely different for each output state, so no mesh is output in this control block.

```
number of surface vars output:
first surface variable ID:
last surface variable ID:
last surface variable ID:
last surface variable ID:
last internal part ID:
last internal part ID:
last user part ID:
la
```

In this domain, the variables are defined at the nodes.

```
solver and domain ID:
number of output vars:
first variable ID:
last variable ID:
last variable ID:
number of DUALCESE3D parts
first internal part ID:
last internal part ID:
partID ndualcese_parts
first user part ID:
last user part ID:
size of each variable component:
number of nodes:
number of nodes:
number of elements:
user partID ndualcese
number of nodal coordinates:
user node numbers:
numelh_dualcese
numelh_dualcese
array of nodal coordinates:
element connectivity:
user element number:
for the first DUALCESE3D element
...
user element number:
```

In this domain, the variables are defined at the surface nodes (of segment sets).

first internal part ID:

last internal part ID:

partID ndualcese_surfsetIDs

first user part ID:

user_surfpartID 1

...

last user part ID:

size of each variable component:

nsurf_dualcese
number of nodes:

number of segments:

user_surfpartID ndualcese_surfsetIDs

nsurf_dualcese

number of nodes:

numelsurf_dualcese

user node numbers:

array of nodal coordinates:

segment connectivity:

partID 1

user_surfpartID ndualcese_surfsetIDs

nsurf_dualcese

numelsurf_dualcese

x_dualcese(nsurf_dualcese)

ix4_dualcese(5, numelsurf_dualcese)

In this domain, the variables are defined at the nodes.

solver and domain ID:

number of output vars:
first variable ID:
last variable ID:
number of DUALCESE parts
first internal part ID:
last internal part ID:
partID ndualcese_parts
first user part ID:
last user part ID:
size of each variable component:
number of nodes:
number of nodes:
number of elements:
user node numbers:
array of nodal coordinates:
element connectivity:
user lement number:
for the first 2D DUALCESE element

first variable ID:
nv_dualcese
ndualcese_parts
ndualcese_parts
numalcese_partID ndualcese_parts
numnp_dualcese
numnp_dualcese
numnp_dualcese
numnp_dualcese
numel2d_dualcese
array of nodal coordinates:
x_dualcese(3, numnp_dualcese)
element connectivity:
for the first 2D DUALCESE element
...

user element number:

In this domain, the variables are defined at the surface nodes (of segment sets).

solver and domain ID:
number of output vars:
first variable ID:
last variable ID:
last variable ID:
last variable ID:
last internal part ID:
last internal part ID:
last internal part ID:
last user p

```
array of nodal coordinates: x_dualcese(3, nsurf_dualcese)
segment connectivity: ix2_dualcese(3, numsurf2Dele_dualcese)
```

In this domain, the variables are defined at the nodes.

```
solver and domain ID:

number of output vars:
first variable ID:
last variable ID:
last variable ID:
last internal part ID:
last internal part ID:
last user part ID ndualcese parts
last user part ID:
las
```

In this domain, the variables are defined at the surface nodes (of segment sets).

```
solver and domain ID:

number of output vars:
first variable ID:
last variable ID:
number of DUALCESE surf. setIDs
first internal part ID:
last internal part ID:
partID ndualcese_surfsetIDs
first user part ID:
user_surfpartID 1
...
last user part ID:
user_surfpartID ndualcese_surfsetIDs
size of each variable component:
number of nodes:
number of segments:
number of segments:
user_surf_dualcese
number of nodal coordinates:
x_dualcese(3, nsurf_dualcese)
segment connectivity:

DUALCESE2DAXI_SURFACE
nv_dualcese

nv_dualcese

ndualcese_surfsetIDs
suser_surfpartID ndualcese_surfsetIDs
nsurf_dualcese
nsurf_dualcese
numsurfel2daxi_dualcese
x_dualcese(3, nsurf_dualcese)
segment connectivity:
ix2_dualcese(3, numsurfel2daxi_dualcese)
```

In this domain, the variables are defined at the particle positions.

General notes regarding "EXTRA DATA TYPES":

The variable IDs are grouped into three groups:

- 1) D3PL_FIRST_SCALAR_ID <= ID < D3PL_FIRST_VECTOR_ID are scalar variables
- 2) D3PL_FIRST_VECTOR_ID <= ID < D3PL_FIRST_TENSOR_ID are vector variables (3 components per entry)
- 3) D3PL_FIRST_TENSOR_ID <= ID < D3PL_END_IDS are symmetric tensor variables (6 component per entry)

When a number of integration points are specified, it is assumed that they are distributed at the Gauss points of the given element type based upon how many are output. That is, for shell or face elements, 4 output points would imply the 2x2 Gauss points are used, while 9 output points would imply the 3x3 Gauss points are used, and so forth. Similarly, for volume elements, 8 output points would imply the 2x2x2 Gauss points are used, while 27 output points would imply he 3x3x3 Gauss points are used, and so forth.

STATE DATA – d3plot and d3part

The state data has three parts:

- Time word and global data
- Node data
- Element data for solids, shell, and beams, respectively

The data defined below can be removed by the flags set in

*DATABASE_EXTENT_BINARY_COMP, see control word INUM =-1 and =-11.

VALUE	LENGTH	DESCRIPTION	
TIME	1	Time word	
GLOBAL	NGLBV	Global variables for this state. LS-DYNA Global Variables: KE, IE, TE, X, Y, and Z velocity IE for each material KE for each material X, Y, and Z velocity for part 1 X, Y, and Z velocity for part n Mass for each part Hourglass energy for each part Force for each rigid wall Xyz position of wall (ls971) = 6 + 7 * (NUMMAT8 + NUMMAT2 + NUMMAT4 + NUMMATT+NUMRBS) + N *	
NODEDATA	NND	NUMRW, N=1, for ls-dyna(ls971) N=4 Total nodal values for state. For LS-DYNA3D and LS-DYNA IT=1, node temperatures only, N=0 IT=2, node temperature and node flux, N=2 IT=3, 3 temperature per node and node flux, N=3 Temperature for shell node at inner, middle and outer layer, inner array, middle array, outer array. IT/10=1, mass scaling value at node. N+=1 =((IT+N)+NDIM*(IU+IV+IA))*NUMNP where IT=temperature flag, IU=coordinates flag, IV=velocities flag, and IA=accelerations flag.	
THERMDATA		If NT3D > 0, NT3D * NEL8 of thermal element data.	
CFDDATA	CFD	Bit flag: NCFDV1, bits from right to left eg, Pressure, Resultant Vorticity, and Density NCFDV1=2+32+1024=1058 2 Pressure 3 X Vorticity	

- 4 Y Vorticity
- 5 Z Vorticity
- 6 Resultant Vorticity
- 7 Enstrophy
- 8 Helicity
- 9 Stream Function
- 10 Enthalpy
- 11 Density
- 12 Turbulent KE
- 13 Dissipation
- 14-20 Eddy Viscosity

Bit flag: NCFDV2

2-11 Species 1 through 10

NOTE: This CFDDATA is no longer output by ls-dyna.

ELEMDATA ENN Total element data for state.

=NEL8*NV3D+NELT*NV3DT+NEL2*NV1D+ NEL4*NV2D+NMSPH*NUM_SPH_VARS The organization of the element data for each element type is described below. The data for the solid elements (7 values/element) is printed first, followed by the data for the beam elements (6 values/element), and then the data for the shell elements (typical 33 or 45 values/element depending on whether the strains are included).

This state data is repeated for each state in the database.

Element data is defined at the integration points within the element. Contour and fringe plots require that the data be extrapolated to or averaged at the nodal points. In LS_PREPOST the element values are averaged at the nodes or optionally extrapolated to the nodes. Element strains are not output by default, these are only output for solids, shell, and thick shell when *DATABASE_EXTENT_BINARY, STRFLG=1

SOLID ELEMENTS – 8 node Hexahedron, other solid elements like wedge, pyramid, and tetrahedron are identified by repeated final connectivities. Eg: pentahedron = 1,2,3,4,5,5,6,6, and tetrahedron = 1,2,3,4,4,4,4.

The database for solid elements consists of 7+NEIPH values per element. NEIPH extra values are defined if and only if NEIPH is greater than zero or if the model is an ALE analysis. If strain components are output, then the last 6 neiph values are true strains: ex, ey, ez, exy, eyz, exz, in the global system.

HIGHER ORDER SOLID ELEMENT STATE DATA

is output after the standard 8 node data. This is only output if either QUADR or CUBIC is > 1. The data is output for each integration point so for example there will be 64 times the stress tensor and plastic strain if the elements used are 64 node hexahedron solids. When this data is output the normal center point data is also output. The data is output according to the order defined in the Higher Order Element data part list.

They are:

- 1. Sigma-x (true stress in the global system)
- 2. Sigma-y
- 3. Sigma-z
- 4. Sigma-xy
- 5. Sigma-yz
- 6. Sigma-zx
- 7. Effective plastic strain or material dependent variable
- 8. First extra value (if NEIPH>0)
- 9. Second extra value (if NEIPH >1)
- 10. Etc. until NEIPH extra values are defined if ISTRN=1
- 7+NEIPH-5. Epsilon-x
- 7+NEIPH-4. Epsilon-y
- 7+NEIPH-3. Epsilon-z
- 7+NEIPH-2. Epsilon-xy
- 7+NEIPH-1. Epsilon-yz
- 7+NEIPH. Epsilon-zx

For thick shell elements the database contains NV3DT = MAXINT* (6 * IOSHL(1) + IOSHL(2) + NEIPS) +12 * ISTRN values per element. Three sets of global stresses are always put into the database for each thick shell and are located at the mid surface, the inner integration point surface, and the outer integration point surface, respectively. If one integration point is used the same through the thickness stress state is outputted three times. If two integration points are used then the mid surface value is taken as the average value. The inner values of the stress are always set to the values at the innermost integration point and likewise for outer values. If the integration point does not lie at the center, ie, an even number of integration points through the thickness, a value is computed that is an average of the two integration point nearest the mid surface.

The IOSHL flags indicate which shell element data is included which is suppressed. The flags are set in ls-dyna by *DATABASE_EXTENT_BINARY, SIGFLG, EPSFLG, RLFLG, and ENGFLG

The ordering of the data follows:

- 1. Sigma-x (mid surface true stress in global system)
- 2. Sigma-y
- 3. Sigma-z
- 4. Sigma-xy
- 5. Sigma-yz
- 6. Sigma-zx
- 7. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for mid surface
- 8. Sigma-x (inner surface true stress in global system)
- 9. Sigma-y
- 10. Sigma-z
- 11. Sigma-xy
- 12. Sigma-yz
- 13. Sigma-zx
- 14. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for inner surface
- 15. Sigma-x (outer surface true stress in global system)
- 16. Sigma-y
- 17. Sigma-z
- 18. Sigma-xy
- 19. Sigma-yz
- 20. Sigma-zx

- 21. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for outer surface
- *. If ISTRN=1, then define strain components Epsilon (x, y, z, xy, yz, zx) here for inner surface and outer surface

If MAXINT > 3 then define an additional (MAXINT-3)*(6*IOSHL(1)+1*IOSHL(2)+NEIPS) quantities here.

For beam elements the database contains NV1D=6 values per element. They are:

- 1. Axial force
- 2. S shear resultant
- 3. T shear resultant
- 4. S bending moment
- 5. T bending moment
- 6. Torsional resultant

If there are values output at beam integration points, then NV1D = 6 + 5 * BEAMIP

- 1. RS shear stress
- 2. TR shear stress
- 3. Axial stress
- 4. Plastic strain
- 5. Axial strain

BEAMIP is set in *DATABASE EXTENT BINARY

If beam material history variables are requested, NEIPB*(3+BEAMIP) number of values are written after the previous data.

Average, min and max, computed from all integration points by LS-DYNA, for NEIPB history variables are written. Then followed by NEIPB history variables for BEAMIP number of integration points.

For example, if NEIPB=2 and BEAMIP=3 the data comes in this order: {hv1_avg, hv2, avg}, {hv1_min, hv2_min}, {hv1_max, hv2_max}, {hv1_ip1, hv2_ip1}, {hv1_ip2, hv2_ip2}, {hv1_ip3, hv2_ip3}.

For shell elements the database contains NV2D values, where:

NV2D=MAXINT* (6*IOSHL(1) + 1*IOSHL(2) + NEIPS) +8*IOSHL(3) + 4*IOSHL(4) + 12*ISTRN values per deformable element. If MATTYP=1 and IRBTYP(I)=20, where I=internal element number, then the material is <u>rigid</u> and the compressed database contains no data for the element. If the minimum value of MAXINT is used, i.e., =3, the stresses are typically located at the mid surface, the inner surface, and the outer surface, respectively. If one integration point is used the stress is written three times. If two integration points are used then the mid surface value is taken as the average value. The inner values of the stress are always set to the values at the innermost integration point and likewise for outer values. If no integration point lies at the center, i.e., an even number of integration points through the thickness, a value is computed that is an average of the two integration point lying nearest the mid surface.

The ordering of the data follows:

- 1. Sigma-x (mid surface true stress in global system)
- 2. Sigma-y
- 3. Sigma-z
- 4. Sigma-x
- 5. Sigma-yz
- 6. Sigma-zx
- 7. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for mid surface
- 8. Sigma-x (inner surface true stress in global system)
- 9. Sigma-y
- 10. Sigma-z
- 11. Sigma-xy
- 12. Sigma-yz
- 13. Sigma-zx
- 14. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for inner surface

- 15. Sigma-x (outer surface true stress in global system)
- 16. Sigma-y
- 17. Sigma-z
- 18. Sigma-xy
- 19. Sigma-yz
- 20. Sigma-zx
- 21. Effective plastic strain or material dependent variable
- *. **Define NEIPS additional history values here for outer surface** If *MAXINT* >3 then define an additional (*MAXINT*-3)* (6*IOSHL(1) + 1*IOSHL(2) + 8*IOSHL(3) + 4*IOSHL(4) + NEIPS) quantities here
 - 22. Bending moment-Mx (local shell coordinate system)
 - 23. Bending moment-My
 - 24. Bending moment-Mxy
 - 25. Shear resultant-Qx
 - 26. Shear resultant-Qy
 - 27. Normal resultant-Nx
 - 28. Normal resultant-Ny
 - 29. Normal resultant-Nxy
 - 30. Thickness
 - 31. Element dependent variable
 - 32. Element dependent variable
 - 33. Internal energy (if and only if ISTRN=0)

The following quantities are expected if and only if ISTRN=1

- 33. eps-x (inner surface strain in global system)
- 34. eps-y
- 35. eps-z
- 36. eps-xy
- 37. eps-yz
- 38. eps-zx
- 39. eps-x (outer surface strain in global system)
- 40. eps-y
- 41. eps-z
- 42. eps-xy
- 43. eps-yz
- 44. eps-zx
- 45. Internal energy (if and only if NV2D>=45)

ELEMENT DELETION OPTION

Skip this section if the word MAXINT is greater than or equal to zero, (MDLOPT>=0). If MDLOPT=1, then the list is equal to the number of nodal points (NUMNP) and contains a one if the node is visible and a zero if the node is not visible, (only used in vec-dyna3d). If MDLOPT=2, then the list equals the total number of elements (NEL8 + NELT + NEL4 + NEL2), in this order, and each value is set to the element material number or =0, if the element is deleted. All these numbers are output as floating point values and not integers.

SMOOTH PARTICLE HYDRODYNNAMICS NODE/ELEMENT STATE DATA

This section is only output if NMSPH>0

For each SPH node the follow values are output:

NUM_SPH_DATA = $1 + \sum isphfg(i)$, i=2:10

Length of data = NUM_SPH_DATA * NUMSPH

Material number, if <=0 then element is deleted.

Currently isphfg(1) = 10, ie number of sph data flags, this could be changed in the future.

If isphfg(2) = 1, - radius of particle influence

If isphfg(3) = 1, - pressure in particle

If isphfg(4) =6, - stress components for particle, sx, sy, sz, sxy, syz, sxz

If isphfg(5) = 1, - plastic strain for particle

If isphfg(6) = 1, - density of particle material

If isphfg(7) = 1, - internal energy of particle

If isphfg(8) = 1, - number of particle neighbors

If isphfg(9) =12, - 6 strain (ex, ey, ez, exy, eyz, exz) and 6 strain rate (erx, ery, erz, erxy, eryz, erxz)

if negative, true strain components

.

If isphfg(10)=1, mass of element (ls971)

Note: it is possible a SPH element could be deleted, or be none active in the initial states, and become active in later states.

PARTICLE STATE DATA (NPEFG > 0)

DES DATA – see description below for state data

STATE DATA

NPARTGAS blocks of NSTGEOM data to describe the state geometry for each bag:

- 1. number of active particles
- 2. current bag volume

PARTICLE DATA

NVAR words of data output for each particle:

- 1. gas ID
- 2. chamber ID
- 3. leakage flag, 0 active, -1 fabric, -2 vent hole, -3 mistracked
- 4. mass
- 5. radius
- 6. spin energy
- 7. translational energy
- 8. distance from particle to nearest segment
- 9. x position
- 10. y position
- 11. z position
- 12. x velocity
- 13. y velocity
- 14. z velocity

ROAD SURFACE MOTION

If NDIM > 5 output rigid body displacement, dx, dy, dz and velocity, vx, vy, vz of each road surface.

Length of data = 6 * NSURF

RIGID BODY MOTION DATA, NDIM=8,9

see DCOMP=5,6 in *DATABASE_EXTENT_BINARY

For each rigid body:

X, Y, Z position of geometric center

MXYZ, Rotation matrix from of principal axes, 3 direction cosines, 9 (3x3) values

VX, VY, VZ translational velocity of GC

RVX, RVY, RVZ, rotational velocity of GC

AX, AY, AZ, translational acceleration of GC

RAX, RAY, RAZ, rotational acceleration of GC

If NDIM=9 only X, Y, Z and MXYZ are output.

EXTRA DATA (Multi-Solver Analysis)

If NCFDV1 = 67108864, then the state data includes NCFDV2 additional datasets from solver-mesh combinations specified after the "User material, node, and element identification numbers" for the structural mesh.

State data of the first solver-mesh combination

...

State data of the last (NCFDV2-th) solver-mesh combination

When the state data comes from the PFEM_IF domain, then the mesh is output first, followed by the data. Currently, the mesh is entirely tetrahedral, but we anticipate users will also specify mixed meshes in the near future:

size of each volume variable component: nnpvol_pfem number of volume nodes: nnpvol_pfem number of tetrahedral elements: ntet_pfem number of pyramid elements: npyr_pfem number of wedge elements: nwdg_pfem number of hexahedral elements: nhex_pfem

user volume node numbers:
array of volume nodal coordinates:
tetrahedral element connectivity:
pyramid element connectivity:
wedge element connectivity:
hexahedral element connectivity:

uvolnodes_pfem(nnpvol_pfem)
xvol_pfem(3, nnpvol_pfem)
ix4_pfem(5, ntet_pfem)
ix5_pfem(6, npyr_pfem)
ix6_pfem(7, nwdg_pfem)
ix8_pfem(9, nhex_pfem)

data for 1st volume variable (size is nnpvol_pfem)

. . .

data for nvolvar_pfem-th volume variable (size is nnpvol_pfem)

Notes:

- (1) the first four entries of ix4_pfem(5, ntet_pfem) for each element are the volume node numbers, while the last entry is the volume part number.
- (2) the first five entries of ix5_pfem(6, npyr_pfem) for each element are the volume node numbers, while the last entry is the volume part number.
- (3) the first six entries of ix6_pfem(7, nwdg_pfem) for each element are the volume node numbers, while the last entry is the volume part number.
- (4) the first eight entries of ix8_pfem(9, nhex_pfem) for each element are the volume node numbers, while the last entry is the volume part number.

When the state data comes from the PFEM_IF_SURFACE domain, then the surface mesh is output first, followed by the data. Currently, the surface mesh is entirely triangular, but we anticipate users will also specify mixed triangle-quadrilateral meshes in the near future:

size of each surface variable component: nnpsurf_pfem

number of surface nodes: nnpsurf_pfem number of surface elements: nelsurf_pfem

user surface node numbers: surfnodes_pfem(nnpsurf_pfem) surface element connectivity: ixsurf_pfem(5, nelsurf_pfem)

data for 1st surface variable (size is nnpsurf_pfem)

...

data for nsurfvar_pfem-th surface variable (size is nnpsurf_pfem)

Notes:

(1) the first four entries of ixsurf_pfem(5, nelsurf_pfem) for each surface element are the surface node numbers, while the last entry is the surface part number. The 3rd and 4th node numbers are the same for triangles.

When the state data comes from the CESE domain, then the mesh is output first, followed by the data:

size of each volume variable component: nele_cese
number of volume nodes: nnpvol_cese
number of tetrahedral elements: ntet_cese
number of pyramid elements: npyr_cese
number of wedge elements: nwdg_cese
number of hexahedral elements: nhex cese

user volume node numbers: volnodes_cese(nnpvol_cese) array of volume nodal coordinates: xvol_cese(3, nnpvol_cese)

tetrahedral element connectivity: ix4_cese(5, ntet_cese)
pyramid element connectivity: ix5_cese(6, npyr_cese)
wedge element connectivity: ix6_cese(7, nwdg_cese)
hexahedral element connectivity: ix8_cese(9, nhex_cese)

data for 1st volume variable (size is nele cese)

...

data for nvolvar_cese-th volume variable (size is nele_cese)

Notes:

- (1) nele_cese = ntet_cese + npyr_cese + nwdg_cese + nhex_cese
- (2) the first four entries of ix4_cese(5, ntet_cese) for each element are the volume node numbers, while the last entry is the volume part number.
- (3) the first five entries of ix5_cese(6, npyr_cese) for each element are the volume node numbers, while the last entry is the volume part number.
- (4) the first six entries of ix6_cese(7, nwdg_cese) for each element are the volume node numbers, while the last entry is the volume part number.
- (5) the first eight entries of ix8_cese(9, nhex_cese) for each element are the volume node numbers, while the last entry is the volume part number.

When the state data comes from the CESE_SURFACE domain, then the surface mesh is output first, followed by the data. The user surface node numbers reference nodes from the CESE domain:

size of each surface variable component: nnpsurf_cese number of surface nodes: nnpsurf_cese number of surface elements: nelsurf_cese

user surface node numbers: surfnodes_cese(nnpsurf_cese) surface element connectivity: ixsurf_cese(5, nelsurf_cese)

data for 1st surface variable (size is nnpsurf_cese)

...

data for nsurfvar_cese-th surface variable (size is nnpsurf_cese)

Notes:

(1) the first four entries of ixsurf_cese(5, nelsurf_cese) for each surface element are the surface node numbers, while the last entry is the surface part number. The 3rd and 4th node numbers are the same for triangles.

When the state data comes from the CESE2D domain, then the mesh is output first, followed by the data:

size of each volume variable component: nele2d_cese number of volume nodes: nnp2d_cese number of triangle elements: ntri_cese number of quadrilateral elements: nquad_cese

user volume node numbers: volnodes_cese(nnp2d_cese)
array of volume nodal coordinates: xvol_cese(3, nnp2d_cese)
triangle element connectivity: ix3_cese(4, ntri_cese)
quadrilateral element connectivity: ix4_cese(5, nquad_cese)

data for 1st volume variable (size is nele2d_cese)

...

data for nvolvar_cese-th volume variable (size is nele2d_cese)

Notes:

- (1) nele2d_cese = ntri_cese + nquad_cese
- (2) the first three entries of ix3_cese(4, ntri_cese) for each element are the volume node numbers, while the last entry is the volume part number.
- (3) the first four entries of ix4_cese(5, nquad_cese) for each element are the volume node numbers, while the last entry is the volume part number.

When the state data comes from the CESE2D_SURFACE domain, then the surface mesh is output first, followed by the data. The user surface node numbers reference nodes from the CESE2D domain:

size of each surface variable component: nnpsurf2d_cese number of surface nodes: nnpsurf2d_cese number of surface elements: nsurfel2d_cese

user surface node numbers: surfnodes_cese(nnpsurf2d_cese) surface element connectivity: ixsurf_cese(3, nsurfel2d_cese)

data for 1st surface variable (size is nnpsurf2d cese)

• • •

data for nsurfvar_cese-th surface variable (size is nnpsurf2d_cese)

Notes:

(1) the first two entries of ixsurf_cese(3, nsurfel2d_cese) for each surface element are the surface node numbers, while the last entry is the surface part number.

When the state data comes from the CESE2DAXI domain, then the mesh is output first, followed by the data:

size of each volume variable component: nele2daxi_cese number of volume nodes: nnp2daxi_cese number of triangle elements: ntri_cese number of quadrilateral elements: nquad_cese

user volume node numbers: volnodes_cese(nnp2daxi_cese) array of volume nodal coordinates: xvol_cese(3, nnp2daxi_cese)

triangle element connectivity: ix3_cese(4, ntri_cese) quadrilateral element connectivity: ix4_cese(5, nquad_cese)

data for 1st volume variable (size is nele2daxi_cese)

...

data for nvolvar cese-th volume variable (size is nele2daxi cese)

Notes:

- (1) nele2daxi_cese = ntri_cese + nquad_cese
- (2) the first three entries of ix3_cese(4, ntri_cese) for each element are the volume node numbers, while the last entry is the volume part number.
- (3) the first four entries of ix4_cese(5, nquad_cese) for each element are the volume node numbers, while the last entry is the volume part number.

When the state data comes from the CESE2DAXI_SURFACE domain, then the surface mesh is output first, followed by the data. The user surface node numbers reference nodes from the CESE2DAXI domain:

size of each surface variable component: nnpsurf2d_cese number of surface nodes: nnpsurf2d_cese number of surface elements: nsurfel2d_cese

user surface node numbers: surfnodes_cese(nnpsurf2d_cese) surface element connectivity: ixsurf_cese(3, nsurfel2d_cese)

data for 1st surface variable (size is nnpsurf2d_cese)

...

data for nsurfvar_cese-th surface variable (size is nnpsurf2d_cese)

Notes

(1) the first two entries of ixsurf_cese(3, nsurfel2d_cese) for each surface element are the surface node numbers, while the last entry is the surface part number.

When the state data comes from the MECH_SURFACE domain, then the surface mesh is output first, followed by the data. The surface node numbers reference nodes from the mechanics mesh:

size of each surface variable component: nnpsurf number of surface nodes: nnpsurf number of surface elements: nelsurf

surface node numbers (internal): surfnodes(nnpsurf) surface element connectivity: ixsurf(5, nelsurf)

data for 1st surface variable (size is nnpsurf)

..

data for nsurfvar-th surface variable (size is nnpsurf)

Notes:

(1) the first four entries of ixsurf(5, nelsurf) for each surface element are the surface node numbers (internal), while the last entry is the surface part number. The 3rd and 4th node numbers are the

same for triangles.

When the state data comes from the MECH2D_SURFACE domain, then the surface mesh is output first, followed by the data. The surface node numbers reference nodes from the mechanics mesh:

size of each surface variable component: nnpsurf2d number of surface nodes: nnpsurf2d number of surface elements: nsurfel2d

surface node numbers (internal): surfnodes(nnpsurf2d) surface element connectivity: ixsurf(3, nsurfel2d)

data for 1st surface variable (size is nnpsurf2d)

...

data for nsurfvar-th surface variable (size is nnpsurf2d)

Notes:

Notes:

(1) the first two entries of ixsurf(3, nsurfel2d) for each surface element are the surface node numbers (internal), while the last entry is the surface part number.

When the state data comes from the DUALCESE3D_ADAPT domain, then the mesh is output first, followed by the data:

```
size of each volume variable component: nnpvol_dualcese
number of volume nodes: nnpvol_dualcese
number of elements: nel_dualcese
user volume node numbers: volnodes_dualcese(nnpvol_dualcese)
array of volume nodal coordinates: xvol_dualcese(3, nnpvol_dualcese)
element connectivity: ix_dualcese(9, nel_dualcese)

data for 1st volume variable (size is nnpvol_dualcese)
...
data for nvolvar_dualcese-th volume variable (size is nnpvol_dualcese)
```

Notes:

(1) the first eight entries of <code>ix_dualcese(9, nel_dualcese)</code> for each element are the volume node numbers, while the last entry is the volume part number. For tetrahedral elements, the last 4 node numbers are the same as the 4th node number. For wedge/prism elements, the 5th and 6th node numbers are the same, and the 7th and 8th node numbers are the same. For pyramid elements, the last 3 node numbers are the same as the 5th node number.

When the state data comes from the DUALCESE_SURF_ADAPT domain, then the surface mesh is output first, followed by the data. The user surface node numbers reference nodes from the DUALCESE ADAPT domain:

```
size of each surface variable component: nnpsurf_dualcese
number of surface nodes: nnpsurf_dualcese
number of surface segments: nelsurf_dualcese
```

(1) the first four entries of ixsurf_dualcese(5, nelsurf_dualcese) for each surface segment are the surface node numbers, while the last entry is the surface setID. The 3rd and 4th node numbers are the same for triangles.

When the state data comes from the DUALCESE2D_ADAPT domain, then the mesh is output first, followed by the data:

```
size of each volume variable component: nnp2d_dualcese
number of volume nodes: nnp2d_dualcese
number of elements: nel2d_dualcese
user volume node numbers:
volnodes_dualcese(nnp2d_dualcese)
array of volume nodal coordinates: xvol_dualcese(3, nnp2d_dualcese)
element connectivity: ix_dualcese(5, nel2d_dualcese)

data for 1st volume variable (size is nnp2d_dualcese)
...
data for nvolvar_dualcese-th volume variable (size is nnp2d_dualcese)
```

Notes:

(1) the first four entries of <code>ix_dualcese(5, nel2d_dualcese)</code> for each element are the volume node numbers, while the last entry is the volume part number. The 3^{rd} and 4^{th} node numbers are the same for triangles.

When the state data comes from the DUALCESE2D_SURF_ADAPT domain, then the surface mesh is output first, followed by the data. The user surface node numbers reference nodes from the DUALCESE2D ADAPT domain:

```
size of each surface variable component: nnpsurf2d_dualcese
                                             nnpsurf2d dualcese
 number of surface nodes:
 number of surface segments:
                                             nsurfel2d dualcese
 user surface node numbers:
surfnodes dualcese(nnpsurf2d dualcese)
 array of nodal coordinates:
                                             x dualcese(3,
nnpsurf2d dualcese)
  surface segment connectivity:
                                             ixsurf dualcese(3,
nsurfel2d dualcese)
  data for 1st surface variable
                                              (size is nnpsurf2d dualcese)
  data for nsurfvar dualcese-th surface variable (size is
nnpsurf2d dualcese)
```

(1) the first two entries of ixsurf_dualcese(3, nsurfel2d_dualcese) for each surface segment are the surface node numbers, while the last entry is the surface setID.

When the state data comes from the DUALCESE2DAXI_ADAPT domain, then the mesh is output first, followed by the data:

```
size of each volume variable component: nnp2daxi_dualcese
number of volume nodes: nnp2daxi_dualcese
number of elements: nel2daxi_dualcese
user volume node numbers:
volnodes_dualcese(nnp2daxi_dualcese)
array of volume nodal coordinates: xvol_dualcese(3, nnp2d_dualcese)
element connectivity: ix_dualcese(5, nel2daxi_dualcese)
data for 1st volume variable (size is nnp2daxi_dualcese)
...
data for nvolvar_dualcese-th volume variable (size is nnp2daxi_dualcese)
```

Notes:

(1) the first four entries of <code>ix_dualcese(5, nel2daxi_dualcese)</code> for each element are the volume node numbers, while the last entry is the volume part number. The 3rd and 4th node numbers are the same for triangles.

When the state data comes from the DUALCESE2DAXI_SURF_ADAPT domain, then the surface mesh is output first, followed by the data. The user surface node numbers reference nodes from the DUALCESE2DAXI ADAPT domain:

```
size of each surface variable component: nnpsurf2d dualcese
 number of surface nodes:
                                             nnpsurf2d dualcese
 number of surface segments:
                                             nsurfel2d dualcese
 user surface node numbers:
surfnodes dualcese(nnpsurf2d dualcese)
                                             x dualcese(3,
  array of nodal coordinates:
nnpsurf2d dualcese)
 surface segment connectivity:
                                             ixsurf dualcese (3,
nsurfel2d dualcese)
  data for 1st surface variable
                                             (size is nnpsurf2d dualcese)
  data for nsurfvar dualcese-th surface variable (size is
nnpsurf2d dualcese)
```

Notes:

(1) the first two entries of ixsurf_dualcese(3, nsurfel2d_dualcese) for each surface segment are the surface node numbers, while the last entry is the surface setID.

When the state data comes from the STOCHASTIC_PARTICLES domain, then the number of particles is output first, following by the particle positions, and then the state data.

```
size of each variable component: n_particles
array of particle positions: x particles(3, n_particles)
```

```
data for 1st output variable
...
data for n prtcl vars-th output variable
```

When the state data comes from the STOCHASTIC PARTICLES domain, then the

```
size of each variable component: n_particles array of particle positions: x_particles(3, n_particles) data for 1st output variable ... data for n_prtcl_vars-th output variable
```

Notes:

There will always be at least the following two variables output for each particle domain: $PARTICLE_SIZES$ and $PARTICLE_VELOCITIES$. That is, $n_prtcl_vars >= 2$. For each particle, both the position and velocity are a 3-component vector.

END OF FILE MARKER

Value = -999999.0 (a floating point number)

TIME HISTORY DATABASE (d3thdt)

There are three sections in the LS-DYNA time history database. The first used to contain 144 words of control information, but now depends upon the number of node and elements the user defines in LS-DYNA. The second contains geometric information including the nodal coordinates and element connectivities. The third section contains the results of the analysis at sequential output intervals for a subset of solids, beams, and shells. The output at a given time, called a state, contains a time word, global variables such as total energies and momenta, nodal data consisting of accelerations, velocities, and displacements, and finally element data is written that may include stresses and strains at integration points. The control information provides information on what is in the file and which database is contained.

CONTROL DATA

VALUE	#WORDS	DISK ADDRESS	DESCRIPTION
Title	10	0	Model Identification
Run time	1	10	Time in seconds since 00:00:00 UTC, January 1, 1970
File type	1	11	1=d3plot, 2=d3drlf, 3=d3thdt, 4=intfor, 5=d3part, 6=blstfor, 7=d3cpm, 8=d3ale, 11=d3eigv, 12=d3mode, 13=d3iter, 21=d3ssd, 22=d3spcm, 23=d3psd, 24=d3rms, 25=d3ftg, 26=d3acs
Source version	1	12	LS-DYNA version *1000000 + svn number
Release number	1	13	Release number in character*4 form 50 for R5.0 511c for R5.1.1c
Version	1	14	Code version, a real number, not integer
NDIM	1	15	Number of dimensions (2 or 3) is set to 4 if element connectivies are unpacked in the LS-DYNA/3D
NUMNP	1	16	Number of nodal points
ICODE	1	17	Flag to identify finite element code =2 old DYNA3D, NIKE3D database =6 new LS-NIKE3D, LS-DYNA/3D database
NGLBV	1	18	Number of global variables to be read in each state
IT	1	19	Flag for temperatures =0 none, =1 read in a temperature for each node

VALUE	#WORDS	DISK ADDRESS	DESCRIPTION
IU	1	20	Flag for current geometry (=1)
IV	1	21	Flag for velocities (=1)
IA	1	22	Flag for accelerations (=1)
NEL8	1	23	Number of 8 node solid elements
NUMMAT8	1	24	Number of materials (parts) used by the 8 node solids
NDS	1	25	Number of node blocks for plotting
NST	1	26	Number of element blocks for plotting. =NSTH + NSTB + NSTS +NSTT
NV3D	1	27	Number of values in database for each solid element
NEL2	1	28	Number of 2 node one-dimensional elements
NUMMAT2	1	29	Number of materials (parts) used by the 2 node 1D elements
NV1D	1	30	Number of values in database for each 1D element
NEL4	1	31	Number of four node two-dimensional elements
NUMMAT4	1	32	Number of materials (parts) used by the 4 node 2D elements
NV2D	1	33	Number of values in database for each 2D element
NEIPH	1	34	Number of additional values per solid element to be written in the type 6 database =NEIPH-6*ISTRN
NEIPS	1	35	Number of additional values per integration point to be written into the type 6 database for shell elements
MAXINT	1	36	Number of integration points dumped for each shell element
NMSPH	1	37	Number of SPH nodes
NGPSPH	1	38	Number of SPH materials
NARBS	1	39	Additional storage required for arbitrary node and element numbering in type 6 database
BLANK	3	40	Unused space
IOSHL(1)	1	43	Stress components flag (=1000 yes)
IOSHL(2)	1	44	Strain components, ISTRN (=1000 yes)
IOSHL(3)	1	45	Shell force resultants (=1000 yes)

VALUE	#WORDS	DISK	DESCRIPTION
		ADDRESS	
IOSHL(4)	1	46	Shell thickness, energy + 2 others (=1000 yes)
BLANK	1	47	Unused space, ignore value
NCFDV1	1	48	Bit flags for CFD nodal values
NCFDV2	1	49	Further bit flags for CFD nodal values
BLANK	7	50	Unused space
EXTRA	1	57	Additional number of control words. If > 0, there are EXTRA control words after the first 64 words.
NSTP	1	58	Number of SPH element blocks
IFLAGD	1	59	Number of node blocks + 1000 flag
NSTH	1	60	Number of solid element blocks
NSTB	1	61	Number of beam element blocks
NSTS	1	62	Number of shell element blocks
NSTT	1	63	Number of thick shell element block
NDSB	2*NDS	64+EXTRA	Node blocks for which time histories are output. The locations 2n-1, where n=1 through NDS correspond to the first node in the block and locations 2n correspond to the last node in the block.
NSTHB	2*NSTH	64+EXTRA+ 2*NDS	Solid element blocks start and end numbers. These are defined in a similar manner to the nodal time history blocks
NSTBB	2*NSTB	64+EXTRA+ 2*NDS+ 2*NSTH	Beam element block start and end numbers
NSTSB	2*NSTS	64+EXTRA+ 2*NDS+ 2*NSTH+ 2*NSTB	Shell element block start and end numbers
NSTTB	2*NSTT	64+EXTRA+ 2*NDS+ 2*NSTH+ 2*NSTB+ 2*NSTS	Thick shell element block start and end numbers
NSTPB	2*NSTP	64+EXTRA+ 2*NDS+ 2*NSTH+ 2*NSTB+ 2*NSTS+ 2*NSTT	SPH element block start and end numbers

ISTRN can only be computed as follows and if NSTS > 0.

If NV2D-MAXINT*(6*IOSHL(1)+IOSHL(2)+NEIPS)+8*IOSHL(3)+4*IOSHL(4) > 10

Then ISTRN = 1, else ISTRN = 0

Or NSTT > 0

If NV3DT-MAXINT*(6*IOSHL(1)+IOSHL(2)+NEIPS) > 10

Then ISTRN = 1, else ISTRN = 0

SMOOTH PARTICLE HYDRODYNAMICS ELEMENT DATA FLAGS

This section is only output if NMSPH > 0. The section is a list of flags to indicate what SPH data is output for each SPH node/element. The first number is the length in words for this array, currently = 10.

SPH elements are centered at nodes, and cover a spherical volume defined by the radius of influence. They do not have a connectivity with other SPH elements. They should be displayed as a dot or a spherical surface, with radius scaling to reduce the size and enable each element to be distinguishable.

As follows:

```
isphfg(1) = 10 - length \ of \ sph \ flags \ array isphfg(2) = 1 - radius \ of \ influence isphfg(3) = 1 - pressure \ in \ particle isphfg(4) = 6 - 6 \ true \ stress \ components isphfg(5) = 1 - plastic \ strain, > 0.0 \ if \ effective \ stress \ exceeds \ yield \ strength isphfg(6) = 1 - density \ of \ particle \ material isphfg(7) = 1 - internal \ energy \ (strain) isphfg(8) = 1 - number \ of \ neighbors \ affecting \ particle isphfg(9) = 12 - 6 \ strain \ (ex, \ ey, \ ez, \ exy, \ eyz, \ exz) \ and \ 6 \ strain \ rate \ (erx, \ ery, \ erzy, \ eryz, \ erzz)
```

if negative, true strain components

```
isphfg(10)=1 - mass of element
```

If the value of isphfg(2-10) = 0, then the particular data item is not output for the particle. To calculated the size of data add the isphfg values from isphfg(2) through isphfg(10) and add one. One value is always output which is the material number as a floating point number for each particle.

If this value is negative then the particle has been deleted from the model.

Full output for each particle is:

```
mat#, radius, pressure, {sx, sy, sz, sxy, syz, sxz} ps, rho, ie, nn, {ex, ey, ez, exy, eyz, exz, erx, ery, erz, erxy, eryz, erxz}, mass.
```

Hence total size is 20.

When a particle is deleted from the model, data is still output for it because the length of data must always be the same for each state.

GEOMETRY DATA

The geometry section contains the nodal coordinates and the element connectivities. The ordering of the nodal points is assumed to be the same as the ordering of the nodal data in the state data that follows. The connectivities are assumed to be packed with 3 integers per word unless NDIM is set to 4 as in the new LS-DYNA/3D, LS-NIKE3D databases. The order of the elements are 3, 2, and 1 dimensional elements if the database is ICODE=2 or 6.

VALUE	LENGTH	DESCRIPTION
X(3,1)	NDIM*NUMNP	Array of nodal coordinates X1,Y1,Z1, X2,Y2,Z2,X3,Y3,Z3,,Xn,Yn,Zn
IX8(9,1)	9*NEL8	Connectivity and material number for each 8 node solid element
IXT(9,1)	9*NELT	Connectivity and material number for each 8 node thick shell element
IX2(6,1)	6*NEL2	Connectivity, orientation node, two null entries, and the material number for each 2 node beam element
IX4(5,1)	5*NEL4	Connectivity and material number for each 4 node shell element

USER MATERIAL, NODE, AND ELEMENT IDENTIFICATION NUMBERS

Skip this section if NARBS (disk address 39) is zero. The user node and element numbers must be in ascending order. *It is assume that if this option is used all the node and element data in the databases is in ascending order in relation to the user numbering.*For sequential material/part numbering, the total length of data is:

NARBS=10+NUMNP+NEL8+NEL2+NEL4+NELT+

3*NMMAT: these numbers are not used

For arbitrary material numbering (NSORT < 0)

NARBS=16+NUMNP+NEL8+NEL2+NEL4+NELT+3*NMMAT

Where material numbers are not in ascending order.

VALUE	LENGTH	DESCRIPTION
NSORT	1	Pointer to arbitrary node numbers in LS-DYNA/3D code, If < 0, it flags that arbitrary material identification numbers are also used.
NSRH	1	Pointer to arbitrary solid element numbers in LS-DYNA code: =NSORT+NUMNP
NSRB	1	Pointer to arbitrary beam element numbers in LS-DYNA code: =NSRH+NEL8
NSRS	1	Pointer to arbitrary shell element numbers in LS-DYNA code: =NSRB+NEL2
NSRT	1	Pointer to arbitrary thick shell element numbers in
		LS-DYNA code: =NSRS+NEL4
NSORTD	1	Number of nodal points
NSRHD	1	Number of 8 node solid elements
NSRBD	1	Number of 2 node beam elements
NSRSD	1	Number of 4 node shell elements
NSRTD	1	Number of 8 node thick shell elements

NSRMA	1	Pointer to an array in the LS-DYNA code that list the material ID's in ascending order.
NSRMU	1	Pointer to an array in the LS-DYNA code that gives the material ID's in the actual order that they are defined in the user input.
NSRMP	1	Pointer to an array in the LS-DYNA code that gives the location of a member in the array originating at NSRMU for each member in the array starting at NSRMA.
NSRTM	1	Total number of materials
NUMRBS	1	Total number of nodal rigid body constraint sets.
NMMAT	1	Total number of materials (parts)
NUSERN	NSORTD	Array of user defined node numbers
NUSERH	NSORTH	Array of user defined solid element numbers
NUSERB	NSORTB	Array of user defined beam element numbers
NUSERS	NSORTS	Array of user defined shell element numbers
NUSERT	NSORTT	Array of user defined solid shell numbers
NORDER	NMMAT	Ordered array of user defined material (parts) ID's
NSRMU	NMMAT	Unordered array of user material (parts) ID's
NSRMP	NMMAT	Cross reference array

TIME HISTORY DATA

The time database contains the following data:

- Time word
- Node data
- Node data for solids, thick shells, and shells, respectively
- Element data for solids, thick shells, beams, and shells, respectively

VALUE	LENGTH	DESCRIPTION
TIME	1	Time word
GLOBAL	NGLBV	Global variables for this state. LS-DYNA Global Variables: KE, IE, TE, X, Y, and Z velocity IE for each part KE for each part X, Y, and Z velocity for part 1 X, Y, and Z velocity for part n Mass for each part Hourglass energy for each part Force for each rigid wall = 6 + 7 * (NUMMAT8+ NUMMAT2 + NUMMAT4 + NUMMATT + NUMRBS) + N*NUMRW, N=1 or N=4 (ls971)

SKIP THE FOLLOWING DATA IF THE NUMBER OF NODE BLOCKS FOR PLOTTING IS ZERO (VALUE NUMDS AT DISK ADDRESS 25)

TIME	1	Time word
NODEDATA	NND	Total nodal values for state where NLN=10*TNODS where TNODS is the number of nodes put into database. The database contains TNODS vectors each with up to 10 components: temperature (if IT=1); x, y, and z coordinates; x, y, and z velocities; and x, y, and z accelerations.

CFDDATA CFD

Bit flag: NCFDV1, bits from right to left Eg Pressure, Resultant Vorticity, and Density NCFDV1=2+32+1024=1058

- 14 Pressure
- 15 X Vorticity
- 16 Y Vorticity
- 17 Z Vorticity
- 18 Resultant Vorticity
- 19 Enstrophy
- 20 Helicity
- 21 Stream Function
- 22 Enthalpy
- 23 Density
- 24 Turbulent KE
- 25 Dissipation
- 14-20 Eddy Viscosity

Bit flag: NCFDV2

2-11 Species 1 through 10

Count number of bits on * NUMNP

SKIP THE FOLLOWING DATA IF THE NUMBER OF ELEMENT BLOCKS FOR IS ZERO (VALUE NUMDS AT DISK ADDRESS 26)

****SKIP THE FOLLOWING IF THERE IS NO DATA FOR SOLID ELEMENTS ****

TIME 1 Time word

SOLIDDATA ENV Total nodal values for solid elements

where ENV=56*TBELM where THELM is the total number of solid elements to be put into the database. The data contains THELM vectors each with 56 components ordered as follows: 8 connectivities: x,y,z coordinates for each of the 8 nodes; and, lastly, x,y,z velocities for each of the 8 nodes.

For solid elements the database contains (7+NEIPH-6*ISTRN) values per element. One set of global stresses are always put into the database for each solid element followed by NEIPH history values. Only data for elements defined in the time history blocks is output. The ordering of the data follows:

- 1. Sigma-x (true stress in the global system)
- 2. Sigma-y
- 3. Sigma-z
- 4. Sigma-xy
- 5. Sigma-yz
- 6. Sigma-zx
- 7. Effective plastic strain or material dependent variable
- 8. First extra value (if NEIPH>0)
- 9. Second extra value (if NEIPH >1)
- 10. Etc. until NEIPH extra values are defined if ISTRN=1
- 7+NEIPH-5. Epsilon-x
- 7+NEIPH-4. Epsilon-y
- 7+NEIPH-3. Epsilon-z
- 7+NEIPH-2. Epsilon-xy
- 7+NEIPH-1. Epsilon-yz
- 7+NEIPH. Epsilon-zx

****SKIP THE FOLLOWING IF THERE IS NO DATA FOR THICK SHELLS ****

TIME 1 Time word

TSHELLDATA ENV

Total nodal values for thick shell elements where ENV=56*TBSEL where TBSEL is the total number of thick shell elements in the database. The data contains TBSEL vectors each with 56 components ordered as follows: 8 connectivities: x,y,z coordinates for each of the 8 nodes; and, lastly, x,y,z velocities for each of the 8 nodes.

****SKIP THE FOLLOWING IF THERE IS NO DATA FOR SHELL ELEMENTS ****

TIME 1 Time word

SHELLDATA ENVS Total nodal values for shell elements where

ENVS=28*TSELM where TSELM is the

total number of shell elements in the database. The data contains TSELM vectors each with 28 components ordered as follows: 4 connectivities: x,y,z coordinates for each of the 4 nodes; and, lastly, x,y,z velocities for each of the 4 nodes.

****SKIP THE FOLLOWING IF THERE IS NO DATA FOR THICK SHELLS ****

For thick shell elements the database contains NV3DT values per element. Three sets of global stresses are always put into the database for each thick shell and are located at the mid surface, the inner surface, and the outer surface, respectively. If one integration point is used the single state is written three times. If two integration points are used then the mid surface value is taken as the average value. The inner values of the stress are always set to the values at the innermost integration point and likewise for outer values. If no integration point lies at the center, i. e. an even number of integration points through the thickness, a value is computed that is an average of the two integration point lying nearest the mid surface. Only data for elements defined in the time history blocks is output. The ordering of the data follows:

- 1. Sigma-x (mid surface true stress in global system)
- 2. Sigma-y
- 3. Sigma-z
- 4. Sigma-xy

- 5. Sigma-yz
- 6. Sigma-zx
- 7. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for midsurface
- 8. Sigma-x (inner surface true stress in global system)
- 9. Sigma-y
- 10. Sigma-z
- 11. Sigma-xy
- 12. Sigma-yz
- 13. Sigma-zx
- 14. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for inner surface
- 15. Sigma-x (outer surface true stress in global system)
- 16. Sigma-y
- 17. Sigma-z
- 18. Sigma-xy
- 19. Sigma-yz
- 20. Sigma-zx
- 21. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for outer surface
- 21. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for outer surface If MAXINT >3 then define an additional (MAXINT-3)* (6*IOSHL(1) + 1*IOSHL(2) + NEIPS) quantities here
 - *. If ISTRN=1, then define strain components Epsilon (x, y, z, xy, yz, zx) here for inner surface and outer surface

****SKIP THE FOLLOWING IF THERE IS NO DATA FOR BEAM ELEMENTS ****

TIME 1 Time word

BEAMDATA BESV Total element values for beam elements

where BESV=NV1D*TBELM. The data contains TBELM vectors each with NV2D

values ordered as described below.

For beam elements the database contains NV1D=6 values per element. They are:

- 1. Axial force
- 2. S shear resultant
- 3. T shear resultant
- 4. S bending moment
- 5. T bending moment
- 6. Torsional resultant

If there are values output at beam integration points, then NV1D = 6 + 5 * BEAMIP

- 7. RS shear stress
- 8. TR shear stress
- 9. Axial stress
- 10. Plastic strain
- 11. Axial strain

BEAMIP is set in *DATABASE EXTENT BINARY

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****SKIP THE FOLLOWING IF THERE IS NO DATA FOR SHELL ELEMENTS ****

For shell elements the database contains NV2D values per element. If the minimum value of MAXINT is 3, then the stresses are typically located at the mid surface, the inner surface, and the outer surface, respectively. If one integration point is used the stress is written three times. If two integration points are used then the mid surface value is taken as the average value. The inner values of the stress are always set to the values at the innermost integration point and likewise for outer values. If no integration point lies at the center, i. e. an even number of integration points through the thickness, a value is computed that is an average of the two integration point lying nearest the mid surface. Only data for elements defined in the time history blocks is output. The ordering of the data follows:

- 1. Sigma-x (mid surface true stress in global system)
- 2. Sigma-y
- 3. Sigma-z
- 4. Sigma-xy
- 5. Sigma-yz
- 6. Sigma-zx
- 7. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for midsurface
- 8. Sigma-x (inner surface true stress in global system)
- 9. Sigma-y
- 10. Sigma-z
- 11. Sigma-xy
- 12. Sigma-yz
- 13. Sigma-zx
- 14. Effective plastic strain or material dependent variable
- *. Define NEIPS additional history values here for inner surface
- 15. Sigma-x (outer surface true stress in global system)
- 16. Sigma-y
- 17. Sigma-z
- 18. Sigma-xy
- 19. Sigma-yz
- 20. Sigma-zx
- 21. Effective plastic strain or material dependent variable
- *. **Define NEIPS additional history values here for outer surface** If MAXINT >3 then define an additional (MAXINT-3)* (6*IOSHL(1) + 1*IOSHL(2) + 8*IOSHL(3) + 4*IOSHL(4) + NEIPS) quantities here

- 22. Bending moment-mx (local shell coordinate system)
- 23. Bending moment-my
- 24. Bending moment-mxy
- 25. Shear resultant-qx
- 26. Shear resultant-qy
- 27. Normal resultant-nx
- 28. Normal resultant-ny
- 29. Normal resultant-nxy
- 30. Thickness
- 31. Element dependent variable
- 32. Element dependent variable
- 33. Internal energy (if and only if ISTRN=0)

The following quantities are expected if and only if ISTRN=1

- 33. eps-x (inner surface strain in global system)
- 34. eps-y
- 35. eps-z
- 36. eps-xy
- 37. eps-yz
- 38. eps-zx
- 39. eps-x (outer surface strain in global system)
- 40. eps-y
- 41. eps-z
- 42. eps-xy
- 43. eps-yz
- 44. eps-zx
- 45. Internal energy (if and only if ISTRN=1)

****SKIP THE FOLLOWING IF THERE IS NO DATA FOR SPH ELEMENTS ****

TIME 1 Time word

Data for each sph element according to the sph flags SPHV=NSTP*NUM_SPH_DATA SPHDATA SPHV

INTERFACE FORCE DATABASE

There are three sections in the interface force database. The first contains 64 words of control information. The second contains geometric information, i.e. the nodal coordinates and segment connectivities for each segment contained in the master and slave surface definitions. The third section contains the results of the analysis at sequential output intervals. The output at a given time is called a state. The state contains a time word, global variables such as total energies and momenta, nodal data consisting of accelerations, velocities, and displacements, and finally segment data is written that include the pressure and shear stress acting on each segment and nodal forces for each node that defines the segment. The control information that follows provides information as to what is in the file and which database is being processed.

CONTROL DATA

			DISK	
VALUE	#W0	ORDS	ADDRES	S DESCRIPTION
Title		10	0	Problem identification
Run time		1	10	time in seconds since 00:00:00 UTC, January 1, 1970
File type		1	11	intfor=4
				1=d3plot, 2=d3drlf, 3=d3thdt, 4=intfor, 5=d3part
				6=blstfor, 7=d3cpm, 8=d3ale, 11=d3eigv,
				12=d3mode, 13=d3iter, 21=d3ssd, 22=d3spcm,
				23=d3psd, 24=d3rms, 25=d3ftg, 26=d3acs
Source versi	on	1	12	ls-dyna version *1000000 + svn number
Release num	nber	1	13	Release number in character*4 form
				50 for R5.0
				511c for R5.1.1c
Version		1	14	Code version
NDIM		1	15	Insert 4 for LS-DYNA/3D database
NUMNP		1	16	Number of nodal points
ICODE		1	17	Insert 6 for LS-DYNA/3D database
NGLBV		1	18	Number of global variable to be read
BLANK		1	19	Insert zero
IU		1	20	Flag for current geometry (=1)

IV	1	21	Flag for velocities (default=1)
BLANK	1	22	Insert zero
BLANK	1	23	Insert zero
BLANK	1	24	Insert zero
BLANK	1	25	Insert zero
BLANK	1	26	Insert zero
BLANK	1	27	Insert zero
BLANK	1	28	Insert zero
BLANK	1	29	Insert zero
BLANK	1	30	Insert zero
NUMSG	1	31	Total number of slave and master segments in sliding interface definitions.
NUMMAT4	1	32	= 2 times the number of sliding interfaces.
NV2D	1	33	= Number of values for each 2D segment.
			If NV2D is negative, then the file is FSIFOR for an
			ALE model
BLANK	5	34	Unused space
NARBS	1	39	Additional storage required for arbitrary node and element numbering in type 6 database This number equals the sum of (10+ NUMNP+NEL8+NEL2+NEL4+ NELT)
BLANK	12	40	Unused space
NHVF	1	52	4*(Number of friction variables)
NTWELD	1	53	4*(Number of tied weld history variables)
NPEN	1	54	0, 12 or 24
NENG	1	55	0 or 4
NTIED	1	56	0 or 2
NWUSR	1	58	4*(Number of wear history variables)
NWEAR	1	59	0, 4 or 8
NPRESU	1	60	0, 1, 2, or 3
NSHEAR	1	61	0 or 3
NFORCE	1	62	0 or 12
NGAPC	1	63	0 or 5

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If *DATABASE_EXTENT_INTFOR is included in the model input the following values apply to the state output, written in the following order:

NV2D = NPRESU + NSHEAR + NFORCE + NGAPC + NWEAR + NWUSR + NHVF +

NTIED + NENG + NPEN + NTWELD

NPRESU: output option for pressures

EQ.0 no pressures output

EQ.1 output normal interface pressure only

EQ.2 output normal interface pressure and peak pressure

EQ.3 output normal interface pressure, peak pressure and time to peak pressure

NSHEAR: output option for maximum interface shear stress,

shear stress in r-direction and s-direction

EQ.0 no

EQ.3 yes

NFORCE: output option for X-, Y- and Z-force at all 4 nodes

EQ.0 no

EQ.12 yes

NGAPC: output option for contact gap at all nodes and frictional energy density

EQ.0 no

EQ.5 yes

NWEAR: output surface wear at each node of segments

EQ.0 no values

EQ.4 output wear depth for each node

EQ.8 output wear depth and sliding distance for each node

NWUSR: Number of wear history variables

4 * Number of wear history values. 4 values for each segment.

NHVF: Number of user friction history variables

4 * Number of user friction history values. 4 values for each segment.

NTIED: Node tied to top face and node tied to bottom face. Value is between 0 (not tied) to 1 (fully tied). May include damage effect.

EQ.0 no values

EQ.2 output tied info for top and bottom side of segment

NENG: Contact energy density. Includes both frictional and contact "springs".

EQ.0 no values

EQ.4 contact energy density for each node

NPEN: Contact penetrations. x,y,z values for each node

EQ.0 no values

EQ.12 Absolute penetrations

EQ.24 Absolute and relative (to max) penetrations

NTWELD: Number of user tied history variables

4 * Number of user friction history values. 4 values for each segment.

GEOMETRY DATA

The geometry section contains the nodal coordinates and the element connectivities. The ordering of the nodal points is assumed to be the same as the ordering of the nodal data in the state data that follows.

VALUE	LENGTH	DESCRIPTION
X(3,1)	NDIM*NUMNP	Array of nodal coordinates X1,Y1,Z1, X2,Y2,Z2,X3,Y3,Z3,,Xn,Yn,Zn
IX4(5,1)	5*NUMSG	Connectivity and identification number for each 3 or 4 node interface segment. For sliding interface n the identification number in 2n-1 for the slave surface and 2n for the master surface.

USER MATERIAL, NODE, AND ELEMENT IDENTIFICATION NUMBERS

Skip this section if NARBS (disk address 39) is zero. The user node and element numbers must be in ascending order. *It is assumed that if this option is used all node and element data anywhere in the databases is in ascending order based on user numbering*. For sequential material/part numbering, the total length of data is:

NARBS=10+NUMNP+NEL8+NEL2+NEL4+NELT+

3*NMMAT: these numbers are not used

For arbitrary material numbering (NSORT < 0)

NARBS=16+NUMNP+NEL8+NEL2+NEL4+NELT+3*NMMAT

Where material numbers are not in ascending order.

VALUE	LENGTH	DESCRIPTION
NSORT	1	Pointer to arbitrary node numbers in LS-DYNA code, If < 0 , it flags that arbitrary material identification numbers are also used.
NSRH	1	Pointer to arbitrary solid element numbers in LS-DYNA code: =NSORT+NUMNP
NSRB	1	Pointer to arbitrary beam element numbers in LS-DYNA code: =NSRH+NEL8
NSRS	1	Pointer to arbitrary shell element numbers in LS-DYNA code: =NSRB+NEL2
NSRT	1	Pointer to arbitrary thick shell element numbers in
		LS-DYNA code: =NSRS+NEL4
NSORTD	1	Number of nodal points
NSRHD	1	Number of 8 node solid elements
NSRBD	1	Number of 2 node beam elements
NSRSD	1	Number of 4 node shell elements
NSRTD	1	Number of 8 node thick shell elements

NSRMA	1	Pointer to an array in the LS-DYNA code that list the contact ID's in ascending order.
NSRMU	1	Pointer to an array in the LS-DYNA code that gives the contact ID's in the actual order that they are defined in the user input.
NSRMP	1	Pointer to an array in the LS-DYNA code that gives the location of a member in the array originating at NSRMU for each member in the array starting at NSRMA.
NSRTM	1	Total number of materials
NUMRBS	1	Total number of nodal rigid body constraint sets
NMMAT	1	Total number of materials (parts)
NUSERN	NSORTD	Array of user defined node numbers
NUSERH	NSORTH	Array of user defined solid element numbers
NUSERB	NSORTB	Array of user defined beam element numbers
NUSERS	NSORTS	Array of user defined shell element numbers
NUSERT	NSORTT	Array of user defined thick shell numbers
NORDER	NMMAT	Ordered array of user defined contact ID's
NSRMU	NMMAT	Unordered array of user contact ID's
Mana		
NSRMP	NMMAT	Cross reference array

ALE, CPM and DES Output Fringe Labels

At the end of the first binary files, eg intfor, fringe labels are appended.

This extra data is written at the end of the following files:

ALE, CPM and DES intfor files. Labels for fringe titles are written directly after the EOF (= -999999.0) marker.

Header output

NTYPE	1	entity type = 90200
NLABEL	1	number of labels
NCHAR8	1	number of char*8 characters in each label

LABEL NCHAR8 Fringe label

For the interface force file (intfor), Fringe labels are written at the end of first file after the EOF (= -999999.0) marker

STATE DATA

The state data for the interface forces have three parts:

- Time word and global data
- Node data
- Force data for sliding interface segments

VALUE	LENGTH	DESCRIPTION
TIME	1	Time word
GLOBAL	NGLBV	Global variables for this state
NODEDATA	NND	Total nodal values for state where NND=(IT+NDIM*(IU+IV))*NUMNP LS-DYNA/3D writes 6 values per node, i.e., the three coordinates and the translational velocities. The Data is put into the database as two vectors: first X(3,NUMNP) and then V(3,NUMNP), respectively.
SEGMDATA	ENN	Data for sliding interface segments where the quantity ENN = NUMSG*NV2D. The organization of the segment data is described below.

This state data is repeated for each state in the database.

For each sliding interface segment the database contains NV2D values per segment. The data order is:

- 1. Normal interface pressure acting on segment
- 2. Maximum interface shear stress acting on segment
- 3. Shear stress in local r-direction of segment
- 4. Shear stress in local s-direction of segment
- 5. X force at node n1 of segment
- 6. Y force at node n1 of segment
- 7. Z force at node n1 of segment
- 8. X force at node n2 of segment
- 9. Y force at node n2 of segment
- 10. Z force at node n2 of segment
- 11. X force at node n3 of segment
- 12. Y force at node n3 of segment

- 13. Z force at node n3 of segment
- 14. X force at node n4 of segment
- 15. Y force at node n4 of segment
- 16. Z force at node n4 of segment
- 17. contact gap at node n1
- 18. contact gap at node n2
- 19. contact gap at node n3
- 20. contact gap at node n4
- 21. surface energy density of segment
- 22. peak pressure
- 23. time to peak pressure
- 24. surface wear depth at node n1
- 25. surface wear depth at node n2
- 26. surface wear depth at node n3
- 27. surface wear depth at node n4
- 28. wear sliding distance at node n1
- 29. wear sliding distance at node n2
- 30. wear sliding distance at node n3
- 31. wear sliding distance at node n4

If contact gap at node n1 = -1.0, then no values are set. Similarly, if surface energy density is = -1.0, no value is set.

If contact gap is >= 999.0 ignore the value, this means that the interfaces are not in discernible contact.

Note: original interface force files have 16 variables, while recent ones, from ls-dyna ls970 and ls971, have up to 31. Check size of NV2D. If *database_extent_intfor is include in the model input then NV2D is according to the formula above.

FSIFOR file output:

- 1. Normal interface pressure acting on segment
- 2. X force on segment
- 3. Y force on segment
- 4. Z force on segment
- 5. relative interface velocity
- 6. X interface velocity
- 7. Y interface velocity
- 8. Z interface velocity

BLSTFOR file (NV2D=16 or 7) output:

- 1. effective (combined incident and reflected) pressure applied to the segment
- 2. reflected wave (relevant only for BLAST=4)
 - =-1: segment is below ground level and not exposed to blast
 - = 0 : segment has not been subjected to blast waves
 - = 1 : segment has been subjected to the initial incident wave
 - = 2 : segment has been subjected to the ground reflected wave
- = 3 : segment resides in the Mach stem region (Since known by geometrical considerations this value is fixed at time t=0. Thus, any non-zero reflected pressure on this segment is due to the Mach wave.)
- 3. incident pressure
- 4. mass density of air
- 5. global x-velocity of blast wind
- 6. global y-velocity of blast wind
- 7. global z-velocity of blast wind

CRACK FILE (d3crck)

The crack file is created in LS-DYNA when the Winfrith Concrete material model is used for solid elements. This model allows up to three orthogonal crack planes to develop each with an origin at the center of the element. The plane is assumed to project to the surface of the solid, and can be represented by drawing a line on any of the six solid element faces where it emerges. This cutting line can be found by considering the intersection of each plane and each solid face.

The crack file is written as a Fortran unformatted binary file, and each record in the file has a start record mark and an end record mark, each of which is 4 bytes for a single precision run and 8 bytes for a double precision run. The data is output for a state and contains:

VALUE	#WORDS	DESCRIPTION
TIME	1	State Time word
NUMBER OF CRACKS	1	NC, Number of sets of crack data
CRACK DATA	16 * NC	Sets of data for each crack

Data for each crack contains: Element ID, Flag for each crack plane, Normal vector for each crack plane and width for each crack plane.

Crack plane flags are as follows:

0 = no crack, 1 = cracked, but no sustaining tensile load, 2 = cracked but closed up, and 3 = fully cracked with no tensile strength.

Shown below is some C coding from LS-PREPOST to illustrate how the crack data is used and the crack lines established. The important coding is picked out in bold type. Also, after the crack data was read in, the crack flags (3 words) where stored as bits in one word. NG and MAT are the Group number for elements and the Material ID we assign in LS-PREPOST.

```
int i, k;
  float x0, y0, z0, x1, y1, z1;
  float xp1, yp1, zp1, xp2, yp2, zp2, pn, dn, t;
  k = 0;
  x0 = xyz[3][0];
  y0 = xyz[3][1];
  z0 = xyz[3][2];
  for (i=0; i<4; i++) {
   /* parametric line clip algorithm */
   xp1 = x0 - xe;
   yp1 = y0 - ye;
   zp1 = z0 - ze;
   x1 = xyz[i][0];
   y1 = xyz[i][1];
   z1 = xyz[i][2];
   xp2 = x1 - x0;
   yp2 = y1 - y0;
   zp2 = z1 - z0;
   x0 = x1;
   y0 = y1;
   z0 = z1;
   pn = nx*xp1 + ny*yp1 + nz*zp1;
   dn = nx*xp2 + ny*yp2 + nz*zp2;
   if (fabs(dn) < 1.0e-15) continue;
   t = -pn / dn;
   if (t < 0.0 || t > 1.0) continue;
   t = t - 1.0;
   xs[k] = x1 + t * xp2;
   ys[k] = y1 + t * yp2;
   zs[k] = z1 + t * zp2;
   k++;
   if (k > 1) break;
  }
  return k;
void SetCrackWidth(float v)
  min_crack_width = MAX(0.0, v);
void DrawCracks(int ist, float *bg_color)
  int i, k, m, n, nc, nd, kd, ip, ic, is;
  int id, facecode;
  int etype, nface, pflag, flag;
  unsigned int j;
  float xc, yc, zc, a, b, c;
  float xyz[24][3];
  float xi[2], yi[2], zi[2];
  int shrink, count, ns[2];
  float dx, dy, dz, ds, d;
  NDCOOR *nod;
  int ng=0;
```

```
float rd, gn, bu;
nod = node;
nod--;
rd = 1.0 - bg \ color[0];
gn = 1.0 - bg_color[1];
bu = 1.0 - bg\_color[2];
glDisable(GL_LIGHTING);
glColor3f(rd, gn, bu);
glLineWidth(2.0);
GetCrackData(ist);
glBegin(GL_LINES);
nc = cstate[ist].nc;
for (n=0; n<nc; n++) {
 pflag = crack[n].pflag;
 if (pflag == 0) continue;
 id = crack[n].id - 1;
 k = solid[id].mat & MASKBIT22;
 j = active_list[k].loc;
 facecode = (active_list[k].akey>>2) & FACEBITS1;
 if (j < BIT30 \&\& facecode > 0) {
  id = j;
  ng = crack[n].ng;
     if (!part[ng].active) continue;
  dx = part[ng].dscale[0];
  dy = part[ng].dscale[1];
  dz = part[ng].dscale[2];
  ds = part[ng].dscale[3];
     shrink = part[ng].shrink_mode;
  etype = (active\_list[k].akey) \& 0x3;
  nface = FACE_NF[etype];
     xc = yc = zc = 0.0; ic = 0;
  for (i=0; i<nface; i++) {
   for (m=0; m<4; m++) {
     nd = FACE_P[etype][i][m];
     kd = solid[id].conn[nd];
        k = 4 * i + m;
        if (ds == 0.0) {
         xyz[k][0] = (disp_state+kd)->xyz[0];
         xyz[k][1] = (disp_state+kd)->xyz[1];
         xyz[k][2] = (disp_state+kd)->xyz[2];
        else {
         GetScaledNodalCoord(kd, nod, disp_state, dx,dy,dz, xyz[k]);
     xc += xyz[k][0];
     yc += xyz[k][1];
     zc += xyz[k][2];
        ic += 1;
    }
     if (ic == 0) continue;
```

```
d = 1.0 / (float)ic; xc *= d; yc *= d; zc *= d;
     if (shrink) {
       for (i=0; i<24; i++) {
        xyz[i][0] = xc + (xyz[i][0] - xc) * shrink_factor;
        xyz[i][1] = yc + (xyz[i][1] - yc) * shrink_factor;
        xyz[i][2] = zc + (xyz[i][2] - zc) * shrink_factor;
  for (i=0; i<nface; i++) {
   if (shrink || (facecode & FACE_CODE1[i])) {
        k = 4 * i;
        /* for each active crack plane */
        for (ip=0,is=0,ic=0; ip<3; ip++,is+=2,ic+=3) {
         flag = (pflag>>is) & 3;
         if (flag == 0) continue;
         if ((min_crack_width < 0.5 && crack[n].cp[ip] >= min_crack_width)
              \|(\min_{\text{crack\_width}} >= 0.5 \&\& \text{ flag} == 3)) 
           a = crack[n].abc[ic];
           b = crack[n].abc[ic+1];
           c = crack[n].abc[ic+2];
           count = CrackPlane(xc, yc, zc, a, b, c, &xyz[k], xi, yi, zi);
           if (count > 1) {
               glVertex3f(xi[0], yi[0], zi[0]);
               glVertex3f(xi[1], yi[1], zi[1]);
          }
glEnd();
glLineWidth(1.0);
```

DYNAIN BINARY FILE FORMAT (dynain.bin)

- /* Discription of Dynain binary format:
- * In first 100 words (integers)
- * head[0] = location of nodal data
- * head[1] = number of nodes
- * head[2] = location of solid element connectivities
- * head[3] = number of solid elements
- * head[4] = location of shell element connectivities + thicknesses
- * head[5] = number of shell elements
- * head[6] = location of adaptive constraints
- * head[7] = number of adaptive constraints
- * head[8] = location of initial stresses for solid elements
- * head[9] = number of initial stress states defined for solids
- * head[10] = location of initial stresses for shell elements
- * head[11] = number of initial stress states defined for shells
- * head[12] = location of initial strains for shell elements
- * head[13] = number of initial strains states defined for shells
- * head[14] = location of boundar spc's
- * head[15] = number of boundary spc's
- * head[16] = location of local coordinate systems by nodes
- * head[17] = number of local coordinate systems by nodes
- * head[18] = location of local coordinate systems by vector
- * head[19] = number of local coordinate systems by vector
- * head[20] = location of initial stress states for beams
- * head[21] = number of initial stress states for beams
- * head[22] = location of thick shell element connectivities
- * head[23] = number of thick shell elements
- * head[24] = location of initial stresses for thick shell elements
- * head[25] = number of initial stress states defined for thick shells
- * head[26] = location of beam element connectivities
- * head [27] = number of beam elements
- * head[28] = location of initial strains for solid elements
- * head[29] = number of initial strain states defined for solids

EXTRA DATA TYPE DEFINITIONS (NCFDV1 = 67108864)

```
#ifndef HAVE D3PLOT
#define HAVE D3PLOT 1
#define D3PL FIRST SCALAR ID 0
#define D3PL FIRST VECTOR ID 1000
#define D3PL FIRST TENSOR ID 2000
#define D3PL END IDS
/* scalar variable names */
enum {
 D3PL Pressure INS=0,
 D3PL Temperature INS,
 D3PL Enstrophy INS,
 D3PL Helicity INS,
 D3PL Stream function INS,
 D3PL Enthalpy INS,
 D3PL Turbulent KE INS,
 D3PL Turbulent eps INS,
 D3PL Eddy Viscosity INS,
  D3PL Density INS,
 D3PL VolFractSpec1 INS,
 D3PL VolFractSpec2 INS,
 D3PL VolFractSpec3 INS,
  D3PL VolFractSpec4 INS,
  D3PL VolFractSpec5 INS,
  D3PL VolFractSpec6 INS,
 D3PL VolFractSpec7 INS,
 D3PL VolFractSpec8 INS,
 D3PL VolFractSpec9 INS,
  D3PL VolFractSpec10 INS,
  D3PL Density CESE,
 D3PL Pressure CESE,
 D3PL Temperature CESE,
 D3PL Total energy CESE,
 D3PL Internal energy CESE,
 D3PL Enthalpy CESE,
  D3PL Entropy CESE,
  D3PL_Stream function CESE,
  D3PL_Density_TS_CESE,
  D3PL Total energy TS CESE,
  D3PL Temperature radflow,
  D3PL Intensity radflow,
  D3PL Scalar potential,
  D3PL Electrical conductivity,
  D3PL Ohm heating power FEM,
  D3PL Ohm heating power BEM,
  D3PL Temperature PFEM,
  D3PL Pressure PFEM,
  D3PL K PFEM,
  D3PL eps PFEM,
  D3PL particle size,
  D3PL particle temperature,
  D3PL particle cnt child particles,
```

```
D3PL Vorticity PFEM,
 D3PL Cp_PFEM,
 D3PL_Qc_PFEM,
 D3PL Shear PFEM,
 D3PL void fraction CESE,
 D3PL schlieren number CESE,
 D3PL LEVELSET PFEM,
 D3PL AVG PRES PFEM,
 D3PL TURB VISC PFEM,
 D3PL relative mu,
 D3PL_HCC PFEM,
 D3PL heatflux PFEM,
 D3PL YPLUS PFEM,
 D3PL UINDEX PFEM,
 D3PL VISCOUS PFEM,
 D3PL RAND r0 EM,
 D3PL_RAND_r10_EM,
 D3PL RAND c10 EM,
 D3PL RAND soc EM,
 D3PL RAND i EM,
 D3PL RAND u EM,
 D3PL RAND v EM,
 D3PL RAND vc EM,
 D3PL_RAND_temperature_EM,
 D3PL RAND P JHR EM,
 D3PL RAND P dudt EM,
 D3PL ALPHA PFEM,
 D3PL CFL PFEM,
 D3PL CMU PFEM,
 D3PL TURBINTENS PFEM,
 D3PL_mass_flow_rate_CESE,
 D3PL Surf Pressure MECH,
 D3PL interface Temperature MECH,
 D3PL Surf solid Heat Flux MECH,
 D3PL Surf fluid Heat Flux MECH,
 D3PL Surf Net HeatFluxRate MECH,
  D3PL HCCAVG PFEM,
 D3PL TemperatureAVG PFEM,
 D3PL HeatfluxAVG PFEM,
 D3PL WETNESS PFEM,
 D3PL RAND areaCircuit EM,
 D3PL RAND areaCell EM,
 D3PL PFEM potential,
 D3PL_Density_DUALCESE,
 D3PL Pressure DUALCESE,
 D3PL Temperature DUALCESE,
 D3PL Tot energy DUALCESE,
 D3PL schlieren no DUALCESE,
 D3PL void fract DUALCESE,
 D3PL Intnl energy DUALCESE,
 D3PL Enthalpy DUALCESE,
 D3PL Entropy DUALCESE,
 D3PL MassFlowRate DUALCESE
/* vector variable names */
enum {
```

```
D3PL Velocity INS=1000,
  D3PL Vorticity INS,
  D3PL Velocity CESE,
  D3PL_Vorticity_CESE,
  D3PL Momentum CESE,
  D3PL Momentum TS CESE,
  D3PL E field radflow,
  D3PL H field radflow,
  D3PL Current density FEM,
  D3PL Electric field FEM,
  D3PL Magnetic field FEM,
  D3PL Lorentz force FEM,
  D3PL Vector potential FEM,
  D3PL_Current_density_BEM,
  D3PL_Electric_field_BEM,
  D3PL_Magnetic_field_BEM,
  D3PL_Lorentz_force_BEM,
  D3PL Vector potential BEM,
  D3PL Surface current,
  D3PL Surface magnetic field,
  D3PL Surface Lorentz force,
 D3PL Velocity PFEM,
 D3PL Vorticity vect PFEM,
  D3PL_particle_velocity,
  D3PL_Average_Velocity_PFEM,
  D3PL H field BEM,
  D3PL magnetization BEM,
  D3PL NF Velocity PFEM,
  D3PL DRAG PFEM,
  D3PL Shear vect PFEM,
  D3PL drag CESE,
  D3PL_Surf_Fluid_Force_MECH,
  D3PL Surf Displacement MECH,
  D3PL Surf Velocity MECH,
  D3PL Surf Acceleration MECH,
  D3PL RAND i vector EM,
  D3PL Shearavg vect PFEM,
  D3PL DRAG element PFEM,
 D3PL Velocity DUALCESE,
 D3PL Vorticity DUALCESE,
 D3PL Momentum DUALCESE,
 D3PL dragforce DUALCESE
};
/* symmetric tensor variable names */
 D3PL INS VELOCITY GRAD=2000
/* chemistry species variable names
                                        */
enum {
 D3PL CHEM SPECIES=3000
```

```
typedef struct d3pnt {
 char * name;
  int id;
} D3PLOT NAME TABLE;
/* Identifiers for solver-mesh combinations */
enum {
 FEM Q1Q0 INS CFD=0,
 CESE CFD NODE,
 CESE CFD ELEMENT,
 CESE CFD ELEMENT TS,
 RADFLOW FULL,
 RADFLOW NODE,
 EM FEMSTER SOLID INTEG PTS,
  EM FEMSTER TSHELL INTEG PTS,
 EM_FEMSTER_SHELL_INTEG_PTS,
 EM FEMSTER SOLID CENTROID,
  EM FEMSTER TSHELL CENTROID,
  EM FEMSTER SHELL CENTROID,
 EM FEMSTER AIR,
 RECT AIR EM NODE,
 EM FEMSTER BEM,
  PFEM IF,
 PFEM IF SURFACE,
 STOCHASTIC PARTICLES,
 CESE SURFACE,
 EM,
  EM SURFACE,
 EM FEMSTER SOLIDSHELL,
 EM FEMSTER NODE,
 CESE2D,
 CESE2D SURFACE,
 CESE2DAXI,
 CESE2DAXI SURFACE,
 CESE SURFACE CFD ELEMENT,
 CESE2D CFD ELEMENT,
 CESE2D SURFACE CFD ELEMENT,
 CESE2DAXI CFD ELEMENT,
 CESE2DAXI SURFACE CFD ELEMENT,
 MECH SURFACE,
 MECH2D SURFACE,
 DUALCESE3D,
 DUALCESE3D SURFACE,
 DUALCESE2D,
 DUALCESE2D SURFACE,
 DUALCESE2DAXI,
 DUALCESE2DAXI SURFACE,
 DUALCESE3D ADAPT,
 DUALCESE3D SURF ADAPT,
 DUALCESE2D ADAPT,
 DUALCESE2D SURF ADAPT,
 DUALCESE2DAXI ADAPT,
 DUALCESE2DAXI SURF ADAPT
};
static D3PLOT NAME TABLE d3plot solver name[] = {
  {"Incompressible FEM CFD", FEM_Q1Q0_INS_CFD},
```

```
{"CESE CFD node", CESE CFD NODE},
  {"CESE CFD element", CESE CFD ELEMENT},
  {"CESE CFD taylor series", CESE CFD ELEMENT TS},
  {"Radiation transport (w/groups)", RADFLOW FULL},
  {"Radiation transport", RADFLOW NODE},
  {"EM solid integ. pts", EM FEMSTER SOLID INTEG PTS},
  {"EM tshell integ. pts", EM FEMSTER TSHELL INTEG PTS},
  {"EM shell integ. pts", EM FEMSTER SHELL INTEG PTS},
  {"EM solid centroid", EM FEMSTER SOLID CENTROID},
  {"EM tshell centroid", EM FEMSTER TSHELL CENTROID},
  {"EM shell centroid", EM_FEMSTER_SHELL_CENTROID},
  {"EM solidShell centroid", EM FEMSTER SOLIDSHELL},
  {"EM node", EM FEMSTER NODE},
  {"EM air", EM FEMSTER AIR},
  {"EM air - rectangular grid", RECT AIR EM NODE},
  {"EM BEM", EM FEMSTER BEM},
  {"Incompressible CFD", PFEM_IF},
  {"Incomp. CFD surfaces", PFEM IF SURFACE},
  {"Stochastic particles", STOCHASTIC PARTICLES},
  {"CESE compressible CFD", CESE},
  {"CESE CFD surface", CESE SURFACE},
  {"EM nodes", EM},
  {"EM surface nodes", EM SURFACE},
  {"CESE 2D CFD", CESE2D},
  {"CESE 2D CFD surface", CESE2D SURFACE},
  {"CESE 2D axisym CFD", CESE2DAXI},
  {"CESE 2D axisym CFD surface", CESE2DAXI SURFACE},
  {"CESE CFD surface element", CESE SURFACE CFD ELEMENT},
  {"CESE 2D CFD element", CESE2D CFD ELEMENT},
  {"CESE 2D CFD surface element", CESE2D SURFACE CFD ELEMENT},
  {"CESE 2D axisym CFD element", CESE2DAXI_CFD_ELEMENT},
  {"CESE CFD axisym surface element", CESE2DAXI SURFACE CFD ELEMENT},
  {"Mechanics surface element", MECH SURFACE},
  {"Mechanics 2D surface element", MECH2D SURFACE},
  {"dual CESE compressible CFD", DUALCESE},
  {"dual CESE surface nodes", DUALCESE SURFACE},
  {"dual CESE 2D compressible CFD", DUALCESE2D},
  {"dual CESE 2D surface nodes", DUALCESE2D SURFACE},
  {"dual CESE 2D axisym comp. CFD", DUALCESE2DAXI},
  {"dual CESE 2D axisym surf nodes", DUALCESE2DAXI SURFACE},
  {"adaptive dual CESE compressible CFD", DUALCESE ADAPT},
  {"adaptive dual CESE surface nodes", DUALCESE SURF ADAPT},
  {"adaptive dual CESE 2D compressible CFD", DUALCESE2D ADAPT},
  {"adaptive dual CESE 2D surf nodes", DUALCESE2D SURF ADAPT},
  {"adaptive dual CESE 2D axisym comp. CFD", DUALCESE2DAXI ADAPT},
  {"adaptive dual CESE 2D axisym surf nodes", DUALCESE2DAXI SURF ADAPT}
static D3PLOT NAME TABLE d3plot et name[] = {
  {"Pressure", D3PL Pressure INS},
  {"Temperature", D3PL Temperature INS},
  {"Enstrophy", D3PL Enstrophy INS},
  {"Helicity", D3PL Helicity INS},
  {"Stream function", D3PL Stream function INS},
  {"Enthalpy", D3PL Enthalpy INS},
  {"Turbulent KE", D3PL Turbulent KE INS},
  {"Turbulent eps", D3PL Turbulent eps INS},
  {"Eddy Viscosity", D3PL Eddy Viscosity INS},
  {"Density", D3PL Density INS},
```

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{"Volume fraction-1", D3PL VolFractSpec1 INS},
{"Volume fraction-2", D3PL VolFractSpec2 INS},
{"Volume fraction-3",D3PL_VolFractSpec3_INS},
{"Volume fraction-4",D3PL_VolFractSpec4_INS},
{"Volume fraction-5",D3PL_VolFractSpec5_INS},
{"Volume fraction-6", D3PL VolFractSpec6 INS},
{"Volume fraction-7", D3PL VolFractSpec7 INS},
{"Volume fraction-8", D3PL VolFractSpec8 INS},
{"Volume fraction-9", D3PL VolFractSpec9 INS},
{"Volume fraction-10", D3PL VolFractSpec10 INS},
{"Fluid velocity", D3PL Velocity INS},
{"Vorticity",D3PL_Vorticity_INS},
{"grad(velocity)",D3PL INS VELOCITY GRAD}
{"Density", D3PL Density CESE},
{"Pressure", D3PL Pressure CESE},
{"Temperature", D3PL_Temperature_CESE},
{"Total energy",D3PL_Total_energy_CESE},
{"Enthalpy", D3PL Enthalpy CESE},
{"Entropy", D3PL Entropy CESE},
{"Stream function", D3PL Stream function CESE},
{"Void fraction", D3PL void fraction CESE},
{"Schlieren number", D3PL schlieren number CESE},
{"Density Taylor series", D3PL Density TS CESE},
{"Total energy Taylor series", D3PL_Total_energy_TS_CESE},
{"Fluid velocity", D3PL Velocity CESE},
{"Vorticity", D3PL Vorticity CESE},
{"Momentum", D3PL Momentum CESE},
{"Momentum Taylor series", D3PL Momentum TS CESE},
{"Temperature radflow", D3PL Temperature radflow},
{"Intensity radflow", D3PL Intensity radflow},
{"E-field radflow",D3PL_E_field_radflow},
{"H-field radflow", D3PL H field radflow},
{"Scalar potential", D3PL Scalar potential},
{"Electrical conductivity", D3PL Electrical conductivity},
{"Ohm heating power FEM", D3PL Ohm heating power FEM},
{"Ohm heating power BEM", D3PL Ohm heating power BEM},
{"Current density FEM", D3PL Current density FEM},
{"Electric field FEM",D3PL Electric_field_FEM},
{"Magnetic field FEM", D3PL Magnetic field FEM},
{"Lorentz force FEM", D3PL Lorentz force FEM},
{"Vector potential FEM", D3PL Vector potential FEM},
{"Current density", D3PL Current density BEM},
{"Electric field", D3PL Electric field BEM},
{"Magnetic field", D3PL_Magnetic_field_BEM},
{"Lorentz force", D3PL Lorentz force BEM},
{"Vector potential", D3PL Vector potential BEM},
{"Surface current", D3PL Surface current},
{"Surface magnetic field", D3PL Surface magnetic field},
{"Surface Lorentz force", D3PL Surface Lorentz force},
{"Fluid velocity", D3PL Velocity PFEM},
{"Fluid temperature", D3PL Temperature PFEM},
{"Fluid pressure", D3PL Pressure PFEM},
{"Fluid vorticity", D3PL Vorticity PFEM},
{"Fluid pressure", D3PL Pressure PFEM},
{"Turbulent K.E.", D3PL K PFEM},
{"Turbulent eps.", D3PL eps PFEM},
{"Particle size", D3PL particle size},
{"Particle velocity", D3PL particle velocity},
{"Particle temperature", D3PL particle temperature},
```

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{"# of child particles",D3PL particle cnt child particles},
{"Pressure Coefficient", D3PL Cp PFEM},
{"Q Criterion", D3PL Qc PFEM},
{"Surface Shear", D3PL Shear PFEM},
{"Level Set", D3PL LEVELSET PFEM},
{"Viscosity", D3PL VISCOUS PFEM},
{"Average Pres", D3PL AVG PRES PFEM},
{"Turbulent Visc", D3PL TURB VISC PFEM},
{"H field", D3PL H field BEM},
{ "Magnetization", D3PL magnetization BEM},
{"Relative permeability", D3PL relative mu},
{"Heat Trans. Coefficient", D3PL HCC PFEM},
{"Heat Flux", D3PL heatflux PFEM},
{"Y plus", D3PL YPLUS PFEM},
{"Uniformity Index", D3PL UINDEX PFEM},
{"Near Field Vel.", D3PL_NF_Velocity_PFEM},
{"Drag Distribution", D3PL_DRAG_PFEM},
{"Surface Shear Vector", D3PL Shear vect PFEM},
{"randle r0", D3PL RAND r0 EM},
{"randle r10", D3PL RAND r10 EM},
{"randle c10", D3PL RAND c10 EM},
{"randle soc", D3PL RAND soc EM},
{"randle i", D3PL RAND i EM},
{"randle u", D3PL_RAND_u_EM},
{"randle v", D3PL RAND v EM},
{"randle vc", D3PL RAND vc EM},
{"randle temp", D3PL RAND temperature EM},
{"randle r0*I^2", D3PL RAND P JHR EM},
{"randle I*T*dudt", D3PL RAND P dudt EM},
{"Thermal diffusivity",D3PL_ALPHA_PFEM},
{"CFL number", D3PL CFL PFEM},
{"RANS Cmu", D3PL CMU PFEM},
{"Turbulent Intens", D3PL TURBINTENS PFEM},
{"Surface mass flow rate", D3PL mass_flow_rate_CESE},
{"Surface drag", D3PL drag CESE},
{"Surface fluid FSI pressure", D3PL Surf Pressure MECH},
{"Surface interface temperature", D3PL interface Temperature MECH},
{"Surface solid displacement", D3PL Surf Displacement MECH},
{"Surface solid velocity", D3PL Surf Velocity MECH},
{"Surface solid acceleration", D3PL Surf Acceleration MECH},
{"Surface fluid FSI force", D3PL Surf Fluid Force MECH},
{"Surface solid heat flux", D3PL Surf solid Heat Flux MECH},
{"Surface fluid heat flux", D3PL Surf fluid Heat Flux MECH},
{"Surface net heat flux rate",D3PL Surf Net HeatFluxRate MECH},
{"Average HTC", D3PL HCCAVG PFEM},
{"Average Temp.", D3PL TemperatureAVG PFEM},
{"Average Heat Flux", D3PL HeatfluxAVG PFEM},
{"randle i vector", D3PL RAND i vector EM},
{"Avg Surface Shear Vector", D3PL Shearavg vect PFEM},
{"Drag Force", D3PL DRAG element PFEM},
{"Wetness", D3PL WETNESS PFEM},
{"Randle areaCircuit", D3PL RAND areaCircuit EM},
{"Riandle areaCell", D3PL RAND areaCell EM},
{"Potential", D3PL PFEM potential},
{"Internal energy", D3PL Internal energy CESE},
{"Density", D3PL Density DUALCESE},
{"Pressure", D3PL Pressure DUALCESE},
{"Temperature", D3PL Temperature DUALCESE},
{"Total energy", D3PL Tot energy DUALCESE},
```

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{"Schlieren_number",D3PL_schlieren_no_DUALCESE},
   {"Void fraction",D3PL_void_fract_DUALCESE},
   {"Internal energy",D3PL_Intnl_energy_DUALCESE},
   {"Enthalpy",D3PL_Enthalpy_DUALCESE},
   {"Entropy",D3PL_Entropy_DUALCESE},
   {"Surface mass flow rate",D3PL_MassFlowRate_DUALCESE},
   {"Fluid_velocity",D3PL_Velocity_DUALCESE},
   {"Vorticity",D3PL_Vorticity_DUALCESE},
   {"Momentum",D3PL_Momentum_DUALCESE},
   {"Surface drag force",D3PL_dragforce_DUALCESE}
};
#endif
```

DES CONTROL BLOCK AND DATA FORMAT

NPEFG - Word #54 : when the seventh digit is set to 1 or 2, there is a DES output. i.e. xx1xxxxxx or V = NPEFG/10000000 = 1 or 2

V = 1: version 1, 8 words in the control block, with words 9 to 16 = 0. (not used)

V = 2: version 2, 16 words in the control block.

1. EXTENDED MASTER CONTROL WORD

There will be one word as the master control word (NDEDB). It defines the number of "Extended Control Blocks" in Section 2, and the same number of "State Data Blocks" in Section 3.

2. DES EXTENDED CONTROL BLOCK

2.1 First 8 words in the block are:

VALUE	WORD#	DESCRIPTION
NDEGP	1	number of DE Parts in Geometries Block
NDEGE	2	number of DE Elements in Geometry Block
NDESP	3	number of DE Parts in State Block
NDESE	4	number of DE Elements in State Block
NGPV	5	number of Part Geometry Variables
NGEV	6	number of Element Geometry Variables
NSPV	7	number of Part State Variables
NSEV	8	number of Element State Variables
version 2:		
BDID	9	database id =0 for Discrete elements
reserved	10-16	not used

2.2 Output Variable Definitions:

N = NGPV + NGPV + NSPV + NSEV

- 2.2.1 There will be N words of output type flags (integers) for each variable listed: The data type flag is in a 6-digit form: GGGCTF, where digits are as following:
 - F: =1, for integers; =2, for real numbers; =3, for complex numbers
 - =5, for array of integers; =6, for array of reals; =7, array of complexes
 - =8, for combined data
 - T: =0, for a scalar variable; =1, for a global vector; =3, for a global tensor
 - =5, for a local vector; =7, for a local tensor; =8, for array data

Notes: A data block contains variable number and types of data, and is included only in the extended control block.

The first integer of a data block (N) is the length of the data block, excluding the first integer. The definition of the data block identified by the group ID.

The post processor may ignore the data block by skipping these (N+1) words, if the definition of the data block is unknown.

- C: The component ID of the grouped variables, defined by group ID,
 - =0: for a scalar variable.
 - =1,2,3: components (X, Y, Z) of a vector variable.
 - =1,2,3,4,5,6: components (XX, YY, ZZ, XY, YZ, ZX) of a symmetric tensor
 - =7,8,9: components (ZY, XZ, YX) of a spin tensor.
 - =1,2,3,4,5,6,7,8,9: components (XX, YY, ZZ, XY, YZ, ZX, ZY, XZ, YX) of a general tensor.

Note: If any components of a vector or tensor not present it is zero.

If none of components (7,8,9) are present, the tensor is symmetrical

If only components (7,8,9) are present, the tensor is unsymmetrical

If components (7,8,9) are present with other components the tensor is general.

GGG: The group ID of the grouped variables (GID)

2.2.2 There will be 8 * N words of the names output for variables listed after the type output (character*8).

2.3 **DES GEOMETRY BLOCK**

2.3.1 There will be NGPV blocks describing geometry data for each DES Part. The block size varies for each part, if variable data blocks are included, i.e. $1+Sum{NGPVi+1, i=2,NGPV}$.

word #1: Part ID

number of words reserved for a DE Part (NGPVi) word #2:

word #3: NGPVi+2: Reserved for a DE element

There will be NGEV blocks describing geometry data for DE Elements. 2.3.2 The block size varies for each element, if variable

data blocks are included, i.e. 5+Sum{NGEVi+1,i=6,NGEV}.

word #1: Element ID
word #2: Part ID
word #3: Radius
word #4: Mass
word #5: Inertia

word #6: number of words for reserved for a DE element (NGEVi)

word #7: NGEVi+6: Reserved for a DE element

3. DES STATE DATA BLOCK

There are no arrays or data blocks in the state data.

3.1 Header for each state: current information for each part: there will be NDESP blocks of part geometry data for each part in. Data output for each part in a state is determined by word #7 (NSPV) from the DES control block. word # 1: Number of active DE elements in each part. to word # NSPV:

3.2 Following each state header: DE Element Data
Data for NDESE DE elements in this state will be listed. Each element will have NSEV (Word #8) words of data output in the database.

The first variables should be always in the database word # 1: flag, 0 active, <0 inactive to word # NSEV:

D3ACS

The binary plot file D3ACS gives the results of frequency domain finite element acoustic analysis. The analysis is activated by the keyword:

*FREQUENCY_DOMAIN_ACOUSTIC_FEM.

The results of acoustic analysis are given as complex acoustic pressure at the nodes. The real and imaginary parts of the pressure, the magnitude of the pressure and a sound pressure level (dB) are written to D3ACS as nodal state variables for each output frequency. No any element state variables are output.

1. Header file

- 1. File type = 26
- 2. There are no element data, values: NV1D, NV2D, NV3D and NV3DT all =0
- 3. Nodal data flags are: IU=1, IT=0, IV=1, and IA=1.

2. State data

The state data is as follows for each frequency:

Time word = Frequency value

NGLBV as given for d3plot

NODEDATA = (IT+NDIM*(IU+IV+IA))*NUMNP

CFDDATA = 0

ELEMDATA = 0

The ordering NODEDATA is as follows:

IU data: X, Y, Z Coordinates for each Node.

IV data: Real part, Imaginary part and Magnitude of Nodal Acoustic Pressure for each Node.

IA data: Sound Pressure Level (dB), Real part and Imaginary part of Nodal Normal Velocity for each Node.