

SHIELD-HIT12A - User's Guide

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- *Поднимите мне веки!*

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- The Institute for Nuclear Research RAS, Moscow, Russia
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Contents

1	Introduction	1
1.1	SHIELD-HIT12A	1
1.1.1	Particle transportation	2
1.1.2	Nuclear physics	2
1.1.3	Atomic physics	3
1.1.4	Scoring of physical quantities	3
1.1.5	Other features	4
2	Quick start guide	5
2.1	Linux: Installation from binary distributions	5
2.2	Windows	5
2.3	Setup a SHIELD-HIT12A run	5
2.4	Example	6
2.5	Analyzing output	9
3	Input files	11
3.1	Overview	11
3.2	<i>mat.dat</i> - Target medium	12
3.2.1	Loading external stopping power files	12
3.2.2	Example of <i>mat.dat</i>	13
3.2.3	List of descriptors for <i>mat.dat</i>	15
3.3	<i>beam.dat</i> - Projectile, statistics, seeds...	16
3.3.1	List of descriptors for <i>beam.dat</i>	17
3.3.2	USECBEAM - External source file	20
3.3.3	USEPARLEV - External parameter file	20
3.3.4	USERIFI - Ripple filter	21
3.4	<i>geo.dat</i> - Geometry	24
3.4.1	Title	24
3.4.2	Bodies	24
3.4.3	Zones	25
3.5	<i>detect.dat</i> - Auxiliary scoring	29
3.5.1	ZONE - Scoring by zone	29
3.5.2	CYL - Cylindrical scoring	30
3.5.3	MSH - Cartesian scoring	31
3.5.4	PLANE - Scoring by 2D plane	32
3.5.5	DZONE, DCYL, DMSH and DPLANE - Differential scoring	33
3.5.6	DMSHZ and DCYLZ - Spectral files in TRiP98 format	34
3.5.7	Available detectors for scoring	36
3.5.8	Scoring examples	37

4	Output files	39
4.1	Auxiliary output files	39
4.2	Management of output with help of special flags	41
5	Auxiliary programs	43
5.1	Helper scripts	43
5.1.1	<i>shield_clean</i> - Remove generated files	43
5.2	Format conversion	43
5.2.1	<i>shield_detect2ascii</i> - Convert binary data to ascii from <i>detect.dat</i>	43
5.2.2	<i>shield2fluka</i> - Convert geo.dat to FLUKA format	43
5.2.3	<i>ddd.py</i> - Convert cylindrical depth dose binning to DDD format	43
5.2.4	<i>shield_dEdx</i> - Prepare an external stopping power table	43
6	Parallelization	45
6.1	Torque	45
6.2	Condor	46
7	User's reference tables	49
7.1	Units	49
7.2	List of particles	50
7.3	Table of nuclear targets with neutron cross sections	51
7.4	List of material ICRU_ID	53
7.5	List of bodies	58
7.6	Input and output files.	60
7.7	List of PARLEV parameters	61
8	Legacy and additional background information	65
8.1	The legacy format of target medium file <i>mat.dat</i>	65
8.2	History	69

List of Tables

7.1	Units	49
7.2	Particles	50
7.3	Nuclear target identifiers	51
7.4	ICRU materials	53
7.5	Input and output files	60
7.6	PARLEV parameters	63

Chapter 1

Introduction

The Monte Carlo particle transport code SHIELD-HIT¹ is designed to precisely simulate therapeutic beams of protons and ions in biological tissue relevant for ion beam cancer therapy. SHIELD-HIT (Heavy Ion Therapy) evolved from the common SHIELD code that models interactions of hadrons and atomic nuclei in complex extended targets in the energy range up to 1 TeV/nucleon.

SHIELD and SHIELD-HIT employ the same models to simulate nuclear interactions, which were developed at JINR, Dubna ² and INR RAS (Moscow) ³. The models are grouped together in the MSDM-generator (Multi Stage Dynamical Model) which allows to simulate a whole nuclear reaction in the exclusive approach. Contrary to the inclusive approach, every generated secondary particle is tracked and processed until it stops, decays, leaves the simulation universe or interacts destructively. Parameters of all secondaries including residual nuclei are retained, thereby fulfilling all conservation laws. Neutron transport below 14.5 MeV is simulated using the 28 group neutron data system ABBN [1] both in SHIELD and SHIELD-HIT.

SHIELD calculates ionization losses of charged hadrons and nuclear fragments according to the Bethe-Bloch equation. The heavy ion version SHIELD-HI contains also ATIMA stopping powers [2]. The "medical" version SHIELD-HIT includes various models and data sets to compute mean ionization loss, fluctuation of ionization losses and multiple Coulomb scattering. In addition, external stopping power tables (including MSTAR and ICRU tables) can be provided using the libdEdx [3] stopping power library.

A more elaborate history of the various releases of SHIELD-HIT is given in section 8.2.

1.1 SHIELD-HIT12A

SHIELD-HIT12A was forked from SHIELD-HIT08 and its source code includes massive changes. The physics engine was tuned to new experimental nucleus-nucleus interaction data, many performance improvements were made and generally the entire code base was restructured, now eliminating the need for users to recompile code for typical usage applications. Several bug fixes from the original SHIELD-HIT branch are forwarded and included in SHIELD-HIT12A. Further development of the Physics engine and bug fixes from the original SHIELD-HIT branch enter continually in SHIELD-HIT12A and vice versa.

Characteristic features of SHIELD-HIT12A can be summarized in the following way:

¹Official homepage: <http://www.inr.ru/shield>

²<http://www.jinr.ru/>

³<http://www.inr.troitsk.ru/>

1.1.1 Particle transportation

- Transport of neutrons, pions, kaons, atomic nuclei, and a number of anti particles in energy range up to 1 TeV/nucleon for SHIELD-HIT. SHIELD-HIT12A is limited to 2 GeV/A. Lower energy cutoff is $E_{\text{cut}} = 25$ keV/nucleon for either version.
- Available initial beam configurations: Gaussian, flat square or flat circular in any direction. Beam divergence can be Gaussian or flat beam; focused or defocused beams can be specified.
- Geometric configuration of the target as an arbitrary combination of geometric bodies bounded by the second order surfaces (Combinatorial Geometry compatible) [4, 5].
- Arbitrary chemical and isotopic composition of materials in target zones can be defined based on the table of available isotopes shown in table 7.3.
- Full memorization of the extra-nuclear cascade tree during simulation without any loss of physics information.
- Formation of neutrons ($E_n < 14.5$ MeV) as well as electrons/positrons and γ -rays, which were created during simulation of extra-nuclear cascade and π^0 decay. However, only neutrons are transported.
- The possibility to switch on/off various physics processes (energy straggling, multiple scattering, nuclear interactions) on user request.

1.1.2 Nuclear physics

- Simulation of inelastic hadron-nucleus and nucleus-nucleus interaction in exclusive approach using a Multi Stage Dynamical Model (MSDM-generator) [6]. The total and inelastic cross sections of the hadron-nucleus and nucleus-nucleus interaction are calculated according to [7–9] with modifications by [10–12]. These cross sections are used for sampling of the nuclear interaction path length as well as for choice of the interaction type (inelastic/elastic). The MSDM-generator describes all stages of the nuclear reaction in the exclusive approach. Current versions of known Russian nuclear models are interfaced in the MSDM-generator:
 - Fast, cascade stage of the nuclear reaction
 - * Intranuclear cascades is handled by the Dubna Cascade Model (DCM) [13]
 - * Independent quark-gluon string model (QGSM) [14–16]
 - * The coalescence model [13]
 - Precompound emission of nucleons and lightest nuclei [17]
 - Equilibrium de-excitation of the residual nucleus
 - * Fermi break-up of light nuclei [18]
 - * Evaporation/Fission competition [18, 19]
 - * Multifragmentation of highly excited nuclei (SMM) [20]
- Neutron transport ($E_n < 14.5$ MeV) on the basis of the 28-group neutron data ABBN [1]. A phase space file of these neutrons can be exported for use in other programs such as MCNP or MCNPX [21].
- Two- and three-particle modes of decay of pions and kaons.

- Neutron transport below 14.5 MeV is simulated by the original neutron transport code LOENT (Low Energy Neutron Transport) [22] using the 28 group neutron data system ABBN [1]. The LOENT code may be used both separately and as a part of the SHIELD code, since SHIELD and LOENT use the same geometry parser (Combinatorial Geometry) [4, 5].

The LOENT code uses the following tables from the neutron data system ABBN:

- st - total cross section
- sf - fission cross section (n,f)
- n - mean number of fission neutrons
- sc - capture cross section (n,c)
- sin - inelastic scattering cross section (n,n'), including the reaction (n,2n)
- se - elastic scattering cross section (n,n)
- m - mean cosine of the angle of the elastic scattering
- sin(g,g+k) - matrix of inter group transitions at the inelastic scattering.

The LOENT code gets neutrons from an external neutron source and follows them, one by one, until the end of the neutron trajectory. Multiplication of neutrons is possible via the reactions (n,2n) and (n,f). Each neutron has its statistical weight attached and a cumulative timer, which accumulates the time from the beginning of the neutron transport history. After transition of the neutron to the thermal group, its energy does not change in further collisions.

1.1.3 Atomic physics

- Ionization losses of charged hadrons and nuclear fragments according to the Bethe-Bloch equation and the Lindhard Scharff equation at low energies.
- Multiple Coulomb scattering simulated with Moliere's or a Gaussian model, similar to that implemented in Geant [23].
- Fluctuations of the ionization energy loss (energy straggling) simulated by a fast implementation (article in preparation) of Vavilov's model [23, 24] or Gaussian distribution.
- Possibility to load external mass stopping power tables to override internal ones. A script interfacing to libdEdx [3] is available.

1.1.4 Scoring of physical quantities

A generic scoring system was developed—that is specified by the user—which provides:

- Detectors for a wide range of physical quantities, which can be made sensitive to specific particles or particle groups.
- Arbitrary scoring in Cartesian or cylindrical scoring grids.
- Alternatively, detectors can be assigned to zones that constitute the target geometry.
- Possibility to generate spectra and depth-dose files in TRiP98 format (*.spc* and *.ddd*).
- Routines to merge results from multiple runs from parallel processing system.

In addition to the generic scoring system, the legacy scoring system is retained, providing:

- Scoring of the production rates of radioisotopes in the entire system.
- Track Length Estimation (TLE) of differential (differential in energy) fluences and energy fluence of secondary particles and nuclear fragments averaged over each geometric zone of the target.
- Scoring of contributions to the energy deposition from various types and from different generations of particles and nuclear fragments separately.

However, the post-processing scripts available for this scoring system cannot merge the ASCII output from parallel runs.

1.1.5 Other features

- Reading and processing of beam source and ripple filter files (see sections 3.3.2 and 3.3.4, respectively).
- Support for parallelization (see chapter 6).

Chapter 2

Quick start guide

2.1 Linux: Installation from binary distributions

SHIELD-HIT12A should come in a tarball of the form: *shield_hitXXa_rXXX.tar.gz*, where *XXX* refers to the SVN revision number. Untar the file:

```
$ tar xvfz shield_hitXXa_rXXX.tar.gz
$ cd shield_hitXXa_rXXX
$ sudo make install
```

which will copy the binaries to */usr/local/bin*. This should work on Ubuntu type systems. If you wish to install to other directories, you may have to edit the Makefile accordingly. On non-sudo type systems (such as Debian or RedHat based Linux systems) one can become root with **su** instead of **sudo** and the last line then reads:

```
$ su
# make install
```

2.2 Windows

SHIELD-HIT12A was tested on Windows 7 and Windows XP systems, but may in principle also run on other Windows systems. SHIELD-HIT12A is provided as a zip file for either 32 or 64 bit systems. Once you unzipped it, you will find a ready-to-run *shieldhit.exe* file.

2.3 Setup a SHIELD-HIT12A run

You can run one of the examples found in base directory, simply by entering:

```
$ shieldhit examples\simple
```

or change into the directory and run **shieldhit** without arguments.

Windows users can open a shell (Run - "Enter a command" - cmd) and change to the directory of where the *shieldhit.exe* file is found. To run the example, simply type:

```
$ shieldhit examples\simple
```

Or simply drop the *shieldhit.exe* file into the folder you want to run, and double click on *shieldhit.exe*.

2.4 Example

Here is a quick description of the most important files in the example *examples/simple* supplied with the distribution:

geo.dat

The file *geo.dat* describes the simulated geometry. Here we want to simulate a target which is a cylinder where the center of its base is located in origin (0,0,0) cm. The height of the cylinder is described by a vector along the Z axis, and is here set to 30 cm. Finally the cylinder radius is set to 10 cm. The cylinder is surrounded by a universe consisting of vacuum which does not interact with the particle. Finally that vacuum is surrounded by a black hole medium. Any particle hitting this medium will be disintegrated, and not tracked any further.

All this information is formatted in “cards”. Cards are lines which are no longer than 80 characters long—a legacy from last centuries punch card systems. Only ASCII characters are allowed; as a white character only space is allowed while tabs are not.

Each card has a number of arguments, which are data being passed to the code. The position and lengths of the arguments varies, thus one should must make sure that the exact positions are used.

Lines beginning with * are ignored, and can be used for comments. So, in the case of the simulation described earlier, our file looks like this:

```
*---><---><-----><----->
      0      0      C12 200 MeV/A, H2O 30 cm cylinder, r=10, 1 zone
*---><---><-----><-----><-----><-----><----->
RCC    1      0.0      0.0      0.0      0.0      0.0      30.0
      10.0
RCC    2      0.0      0.0      -5.0      0.0      0.0      35.0
      15.0
RCC    3      0.0      0.0      -10.0      0.0      0.0      40.0
      20.0
END
001      +1
002      +2      -1
003      +3      -2
END
      1      2      3
      1 1000      0
```

The first line, is a comment line, which is ignored. It is inserted here, to better see where the fields are which are interpreted by SHIELD-HIT12A. The following line specifies a title of the geometry setup. The title is preceded by two integers which are described later in section 3.4. After another comment line, the next two lines describe the geometry of the aforementioned target cylinder. The **RCC** card selects a body (here a Right Circular Cylinder). The body is assigned to a number 1. The x,y,z coordinates of the base follow, along with a vector which spans to the top of the cylinder. Since the cylinder is following the Z axis, all entries are zero, except for the height of the cylinder, i.e. the z-component of the vector. The following card is a continuation card of the **RCC** card, and holds only one value, namely the radius of the cylinder. All values are in cm. A list of default units in SHIELD-HIT12A is provided in section 7.1.

Afterwards two more cylinders follow, each expanded by an additional 10 cm in all three directions.

The geometry section, which only contains one body here, must be terminated by the **END** card.

The lines

```
001      +1
002      +2      -1
003      +3      -2
```

assigns the bodies we just specified to zones. For the first body (our target cylinder) the body and the zone are identical. The surrounding vacuum is represented by the second cylinder, but subtracted the inner cylinder, since this is already described by zone 001. Similarly zone 003 consists of the largest body (number 3), subtracted the body number 2. There is no reason to subtract body number 1 here, as it is already entirely contained in body number 2. The point is that a zone can be any Boolean composition of previously defined bodies, which is described in detail later in section 3.4. Any point inside the black hole must be assigned to a zone, and only one zone.

The section describing the zones must also be terminated with the **END** card.

Finally, two lines follow: the first specifies a list of zones, and the second assigns a material number to that zone. In this example, each line consists only of three numbers since we have three zones. In the second line we tell which materials each zone consists of. The numbers 0 and 1000 are predefined, any other number will be specified in the *mat.dat* file. 0 is black hole, 1000 is vacuum and the inner most cylinder is made of material with the id# 1.

mat.dat

The file *mat.dat* specifies the medium of the zone we just specified in *geo.dat*.

We define medium number 1 to be water. The most simple way to do this is by writing:

```
MEDIUM 1
ICRU 276
END
```

The first line tells SHIELD-HIT12A that we now are going to assign **MEDIUM** number one. The second line says that we want to use a default ICRU material with material number 276. The numbers of predefined materials (as given by ICRU) are listed in table 7.4 in section 7.4 with 276 being water. Each **MEDIUM** section must be terminated with an **END** line for SHIELD-HIT12A to initialize the material. The specification materials, other than those listed in table 7.4, will be explained in section 3.2.

beam.dat

For *beam.dat* we should initially only worry about a few lines. Fairly at the top, **JPART0** specifies the number of the primary particle according to table 7.2.

```
JPART0      25      ! Incident particle type
HIPROJ      12.0      6.0      ! A and Z of heavy ion
```

Here we request a beam of heavy ions (**JPART0** = 25), which is specified further with the **HIPROJ** card. By stating $A = 12$ and $Z = 6$ we select carbon-12 ions to be our primary particle.

The energy is specified with the **TMAX0** card, here set to 391.0 MeV/nucleon.

```
TMAX0      391.0      0.0      ! Incident energy; (MeV/nuc1)
```

The statistics, i.e. the amount of primary particles to be simulated is specified in the `NSTAT` line:

```
NSTAT                2000    1000          ! NSTAT, Step of saving
```

With no further input we get a default pencil beam with $\sigma_x = 0.0$ $\sigma_y = 0.0$ cm which starts at origin (0,0,0).

detect.dat

Finally, we can specify detectors to score a variety of quantities. The file *detect.dat* is optional, but very useful. In this example, we want to score a depth dose-curve along the water target.

We can setup a cylinder, following the Z axis, located inside the water target. It does not need to match any geometry specified in *geo.dat*. Note, this file requires fixed format.

```
*-----0---><-----1---><-----2---><-----3---><-----4---><-----5---><-----6--->
CYL                0.0        0.0        0.0        10.0        7.0        30.0
                   1          1          300         -1        ENERGY    ex_cyl
```

The first card specifies cylindrical scoring, starting from point (0,0,0), and with the radius 10 cm, covering 2π (by specifying “7”¹) of the circumference of the cylinder, and with a height of 30 cm along the Z axis. Currently, the `CYL` scoring cannot be rotated, and is always placed parallel to the X-axis. In the next line the first three arguments say that we want 1 bin for the radius, 1 bin for the angle spanning the cylinder² and 300 bins along the height. The result will then be a one-dimensional file, where the scored quantity is listed in steps of 30.0 cm / 300 bins (i.e. 1 mm). The next three arguments tell that we want the `ENERGY` to be scored of the particle -1 which means the energy deposited by all particles crossing the bins. `ex_cyl` is the name of the output file.

Alternatively we could also specify a cartesian scoring mesh. The two next lines specifies a mesh with one bin from -5 to 5 cm along the X and Y axis, and then 300 bins from 0 cm to 30 cm along Z. The result is saved in the file *ex_zmsh*.

```
MSH                -5.0        -5.0        0.0        5.0        5.0        30.0
                   1          1          300         -1        ENERGY    ex_zmsh
```

We can here also score along another axis if we want. The next example scores 100 bins from -5 cm to 5 cm along the Y axis.

```
MSH                -1.0        -5.0        10.0        1.0        5.0        12.0
                   1          100         1          -1        ENERGY    ex_ymsh
```

or if we want to see a 100x300 pixel 2-D map of the energy deposited along the beam path in the target:

```
MSH                -5.0        -5.0        0.0        5.0        5.0        30.0
                   1          100         300         -1        ENERGY    ex_yzmsh
```

¹Any number larger than 2π means we do not subdivide the cylinder along its radial lines, i.e. we do not cut the cylinder in pieces of pie.

²I.e. scoring along the entire circumference and not into pie pieces.

2.5 Analyzing output

After SHIELD-HIT12A terminated the run you find several new files in your directory. Here the results from the four detectors, defined in *detect.dat*, are stored in the files *ex_zmsg*, *ex_ymsh*, *ex_yzmsb* and *ex_cyl*. These files are in binary form and the script *shield_detect2ascii* can be used to convert them to ASCII format. Stepping into the directory and applying,

```
$ shield_detect2ascii ex_zmsb
```

will produce a human-readable output file *detout.txt*, which can be plotted.

Chapter 3

Input files

3.1 Overview

The SHIELD-HIT12A transport code works with at least three input files.

mat.dat – chemical composition of materials in target zones.

beam.dat – several parameters, explained below (like seed, projectile, statistics, etc.).

geo.dat – geometry of the target using combinational geometry (CG), similar to FLUKA.

Optionally, the user can include an additional file:

detect.dat – for simple scoring of physical quantities in independent geometries.

These four files are described in more detail in the sections to follow.

As extension of the file *beam.dat* up to three optional ASCII formatted files (with user-specified names) can be supplied that specify: a ripple filter, a beam source file, and parameters for nuclear interaction models. Examples for these three files are shown adjacent to the description of file *beam.dat* in 3.3. If chosen by the user, SHIELD-HIT12A may also look for external ASCII files containing stopping power data as described in section 3.2.

All input files must be grouped in the same folder. Its location can be specified by providing a command line argument to the SHIELD-HIT12A executable. All output files are placed into the same directory. The SHIELD-HIT12A executable itself can be run from anywhere if the exact path to the directory with the input files is given. If the executable is run from the directory containing the input files the specification of the path can be omitted.

3.2 *mat.dat* - Target medium

The *mat.dat* file defines the materials that are used in the simulation. It consists of cards (lines starting with a descriptor followed by possible arguments) and comment lines. Any line starting with a * or a < or a blank line are regarded as a comment, which is ignored and may appear at any place. The ordering of the cards has to obey certain rules. Some cards can be omitted, and SHIELD-HIT12A will assign the default values given in the descriptor list below. It is recommended to specify all parameters explicitly since default values may change in future releases of SHIELD-HIT12A. While the ordering of the arguments must be as specified for the respective card, they can be positioned in any column (free format).

Each parseable line starts with a descriptor, which specifies what parameters are to be described. After this descriptor, none, one or two arguments follow, depending on the type of the descriptor. The descriptor and each argument must be separated with at least one space or tab. The rest of the line is not parsed. It can be used for comments if it is separated from the last argument with at least a space or tab. Only the first 128 characters are read; any characters beyond will be ignored.

Each material specification starts with the MEDIUM card. A medium can consist of a single element or of a compound. A compound again can be a chemical compound, or simply a mix of elements and/or isotopes, e.g.:

- Pure medium - element with one fixed Z (e.g. O, Fe, Cu, etc.).
- Chemical composition, e.g. water H_2O .
- Isotope mixture, if isotopes of the same Z have different neutron properties at low energies, e.g. mixture of ^{235}U and ^{238}U .
- a chemical compound containing isotopes, e.g. ^6LiF .

The maximal total number of different media NUMMED is limited to 16, excluding internal and outer vacuum which are predefined as medium numbers 1000 and 0, respectively.

The user can select one of the predefined media listed in the ICRU material table in section 7.4. Alternatively, the user can define a MEDIUM by specifying the chemical elements or isotopes, which can be found in this medium, limited to a total number of 13. A chemical element is described by the variable NUCLID. Usually, $\text{NUCLID}=Z$ selects the element from the periodic table corresponding to the atomic number Z ($1 < Z < 100$). Further isotopes, which differ in the number of neutrons, are predefined with values of $\text{NUCLID} > 100$ and are listed in table 7.3. If an isotope does not appear in the list, no low energy neutron cross section data are included in SHIELD-HIT12A. In this case SHIELD-HIT12A will not start the simulation, if the neutron transport cutoff energy (NEUTRLCUT in *beam.dat* see section 3.3) is set below 14.5 MeV.

Note that the corresponding neutron cross sections are either for natural isotope mixtures or for a specific isotope.

As of SHIELD-HIT12A the *mat.dat* format was changed to the format described above. The old SHIELD-HIT format can still be employed and is described in section 8.1. However, its use is discouraged since this part will not be maintained and may be removed in future. The new format is clearly advantageous since new features can be added without breaking backwards compatibility and it is more convenient to work with. Users still applying the old *mat.dat* format are therefore encouraged to consider migrating to the new file format.

3.2.1 Loading external stopping power files

The user can provide external files containing mass stopping power data by specifying the LOADEDX card with an appropriate value from the material list in table 7.4. Each file con-

tains data for the lightest 18 ions from $^1\text{H}^+$ (proton) to $^{40}\text{Ar}^{18+}$ interacting with one specific target material.

These external files are expected to be ASCII formatted, space or tab separated columns, and should follow the ICRU 73 [25] energy grid (53 energy nodes from 25 keV/u up to 1 GeV/u). SHIELD-HIT12A expects the grid to contain the first 18 ions, i.e. from protons ($Z=1$) to Argon ions ($Z=18$). The mass stopping power data must be in units of $\text{MeV cm}^2/\text{g}$. Comment lines starting with $*$ are allowed anywhere in the file, but no line (either comment or data) may be longer than 512 characters. External stopping power files may contain data different from ICRU, e.g. ATIMA [2] data, but the energy grid must be compatible to ICRU 73. The libdEdx [3, 26] computer library can serve as a source for stopping power data. A script *shield_dEdx*, described in section 5.2.4, can generate these tables for most ICRU materials using libdEdx.

Examples of external stopping power files are: *Water.txt*, *Air.txt*, *A-150.txt*, *Kapton.txt*, which are supplied along with the SHIELD-HIT12A distribution. In table 7.4 the naming convention of the external stopping power files is listed.

3.2.2 Example of *mat.dat*

An example of *mat.dat* with five materials is given below

1. Water, ICRU default setup. External stopping powers will be loaded from *Water.txt*.
2. Air, ICRU default setup.
3. PMMA (also known as Perspex or Lucite), is defined. It is a compound in solid state, here with a density of 1.19 g/cm^3 . The I-value of hydrogen is changed to 21.9 eV.
4. Alanine, ICRU default setup, but overriding the density.
5. Lithium-6 Fluoride (LiF), which is composed of Lithium-6 isotopes which has a very high thermal neutron cross section. External stopping power tables are loaded from *LiF.txt*.

```
* ----- WATER -----
MEDIUM      1
ICRU         276
LOADEDX
END
* ----- AIR -----
MEDIUM      2
ICRU         104
END
* ----- PMMA -----
MEDIUM      3
STATE       2
RHO         1.19
NUCLID      1      8
IVALUE      21.9    ! Override I-value of hydrogen
NUCLID      6      5
NUCLID      8      2
END
* ----- NPL ALANINE -----
MEDIUM      4
ICRU         105
```

```
RHO      1.23      ! Override density, since NPL alanine is less dense
END
* ----- Li6-Fluoride -----
MEDIUM   5
STATE    2
RHO      2.635
NUCLID   105    1
NUCLID   9      1
LOADEDX  185      ! Better stopping powers from external lib.
END
```

3.2.3 List of descriptors for *mat.dat*

The recognized descriptors are listed alphabetically, and not by relevance.

AMASS : 1 argument. Optional card which can override the atomic mass A of an element when specified *after* a **NUCLID** card.

END : No argument. Mandatory card that terminates the description of a medium that was started by the **MEDIUM** card.

ICRU : 1 argument. Optional card which will select a material from the ICRU list in section 7.4. This card may not be used together with the **NUCLID** card for a **MEDIUM**, as it may lead to unpredictable behaviour.

IVALUE : 1 argument. Optional card that can override the default mean excitation energy (I -value) of the element that was specified with the preceding **NUCLID** card. Units are eV.

LOADEDX : 1 optional argument. Optional card which will trigger SHIELD-HIT12A to look for an external stopping power file following the naming scheme stated in the table in section 7.4. The argument must be the ICRU number of the material to be loaded. However, if used after the **ICRU** card, the ICRU number is already known, and the argument can be omitted. The external stopping power files should contain mass stopping power data for the medium for ions ranging from $^1\text{H}^+$ (proton) to $^{40}\text{Ar}^{18+}$. See section 3.2.1 for a description of the format.

MEDIUM : 1 argument which specifies the medium number. The numbering must be sequential, starting at 1 for the first medium which is specified.

NUCLID : 2 arguments. This card adds a single element to the medium. First argument is the **NUCLID** number specifying the element or isotope from table 7.3. You can also use a Z value that is missing in this list but there are no neutron cross section data attached to it. Second argument is the relative stoichiometric fraction, which will automatically be normalized by SHIELD-HIT12A. Up to 13 elements can be defined per **MEDIUM**. SHIELD-HIT12A will refuse to start, if there are no neutron cross sections for the requested element, and the lower neutron energy cut is set below 14.5 MeV.

RHO : 1 argument which specifies the density of the medium in g/cm^3 . This card is mandatory in conjunction with the **NUCLID** card. It should be stated after the **MEDIUM** card and before the first **NUCLID** card. In the case that the material is setup using the **ICRU** card, **RHO** will override the default value.

STATE : 1 arguments which specifies the state of the medium. Appropriate mean excitation energies (I -values) are selected for each element in a compound, following the ICRU49 guidelines [27]. This card is mandatory in conjunction with the use of a **NUCLID** card. It should be stated after the **MEDIUM** card and before the first **NUCLID** card.

SHIELD-HIT12A discriminates between a chemical compound and just a mix of elements. If **STATE** 0, then SHIELD-HIT12A uses I -values for elements. If **STATE** 1 or **STATE** 2 SHIELD-HIT12A uses I -values for compounds in gas or liquid/condensed phase, respectively. In the case that the material is setup using the **ICRU** card, **STATE** will override the default state of the **ICRU** material.

3.3 *beam.dat* - Projectile, statistics, seeds...

The *beam.dat* file characterizes the particle beam transport by defining: the primary beam properties; the required statistics and random seed; details of the physics engine; and additional options.

The format is free, meaning that the arguments can be given in any columns, but they must be ordered as in the description of the respective card. After the last argument, any text string for comments are allowed. Cards can be arbitrarily ordered. Any card (with exception of HIPROJ, see below) can be omitted, and SHIELD-HIT12A will then use default values. The default values are given in the descriptor list below, but may change unnoticed in future releases of SHIELD-HIT12A. Therefore it is recommended to avoid using default parameters, if they are important to the simulation.

In *beam.dat* any line starting with a * or a < is regarded as a comment and is ignored. Blank lines will be ignored too.

Each parsable line starts with a descriptor that specifies what parameters are to be described. After this descriptor, one to three arguments follow, depending on the type of the descriptor. The descriptor and each argument must be separated with at least one space or tab. The rest of the line is not parsed, and can be used for comments. However it is important that the comments are separated too from the last argument with a space or tab. Only the first 128 characters are read, any characters beyond will be ignored.

Example of typical *beam.dat* file:

```
*          Input file beam.dat for the SHIELD-HIT Transport Code
RNDSEED           89736501      ! Random seed
JPARTO            25          ! Incident particle type
HIPROJ            12.0         6.0      ! Carbon ions
TMAXO             400.0        0.0      ! Incident energy; (MeV/nucl)
BEAMSIGMA         2.0          2.0      ! SigmaX, SigmaY at focus point
BEAMPOS           0.0          0.0      -1.50 ! Beam XYZ start pos
NSTAT             20000        5000     ! NSTAT, Step of saving
NEUTRLCUT         0.0          ! Cutoff for neutron transport
WRITEFILE         ynnnnnnnnnnn ! Various flags y=yes, n=no
MAKELN            0           ! 1 - Make neutron phase space file
DELTAЕ            0.020       ! Delta E (relative share ~0.1)
DEMIN             0.030       ! Minimum Energy step 0.025 (MeV/n)
STRAGG            2           ! Straggling: 0-Off 1-Gauss, 2-Vavilov
MSCAT             2           ! Mult. scatt 0-Off 1-Gauss, 2-Moliere
NUCRE             1           ! Nucl.Reac. switcher: 0-Off, 1-On
USERIFI           rifi.dat     ! Filename of ripple filter file
USECBЕAM          sobp.dat     ! Filename of beam sourcefile
```

3.3.1 List of descriptors for *beam.dat*

The recognized descriptors are listed alphabetically, and not by relevance.

BEAMDIR : Optional card. 2 arguments that specify the beam direction: the polar and azimuth angles θ and ϕ , both in degrees, as explained in figure 3.1.

Initially, the beam divergence and shape is setup assuming transport along Z axis. No matter where the card is specified, the actual rotation of the beam into the requested beam direction will always take place as a last step, before transportation begins. If this card is not specified, then beam transport along the Z axis is assumed.

If this card is not specified, θ and ϕ are set to zero. That is, by default the beam points in Z direction.

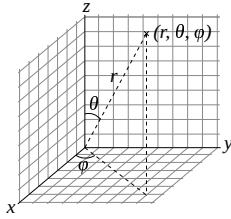


Figure 3.1: Coordinate system as used by SHIELD-HIT12A for describing the beam direction. θ is the angle relative to the Z axis. ϕ is the angle of the orthogonal projection of the direction onto XY plane measured from the X axis.

BEAMDIV : Optional card. 3 arguments that specify the divergence and focal point of the beam. First two arguments define the divergence in Gaussian sigmas in X and Y direction assuming beam transport in Z direction. (A subsequent rotation of the beam can be specified with the **BEAMDIR** card.) The beam divergence must be in mrad. Negative values will generate a flat distribution instead. The distributions are weighted along the polar angle, so requesting a flat angle distribution will also result in a flat beam spot.

The third argument sets the distance to the focal point k in cm at which the undisturbed beam (i.e. transport in vacuum) has the width and shape specified by the **BEAMSIGMA** card. A positive value $k > 0$ describes a defocused beam with the focal point upstream the beam at a distance k , while a negative value $k < 0$ results in a focused beam with the focal point downstream, relative to the source position of the beam.

The beam model follows equation (25) in reference [28], which gives a beam with a slant emittance ellipse. The same reference describes the Gaussian case only, however k affects any beam shape in SHIELD-HIT12A.

If this card is not specified, all values are set by default to zero.

BEAMPOS : Optional card. 3 arguments that specify the start position of the beam in (x,y,z) coordinates, respectively. All units are in cm.

By default the beam starts at the origin (0,0,0).

BEAMSIGMA : Optional card. 2 arguments that specify the lateral extension of a Gaussian-shaped beam by the sigmas in cm along X and Y, respectively. If **both** values are negative, then a flat square distribution is generated, where σ_x and σ_y represent half the width of the sides. If σ_x is larger than 0 **and** σ_y is smaller than 0, then a flat circular distribution is generated with radius $\text{abs}(\sigma_y)$. The σ_x value is unused here, but it must be larger than 0.0. If the focus is specified in the optional **BEAMDIV** card then the beam size and shape are obtained at the focal point.

Default values are zero sigma in X and Y direction.

DELTAE : Optional card. 1 argument that specifies the relative mean energy loss per transportation step, that is a finite fraction smaller than unity. This value times the particle energy at a given step determines the absolute energy loss.

Default is 0.05. (Meaning 5%.)

DEMIN : Optional card. 1 argument that specifies the minimum allowed energy loss per transportation step in MeV/n.

Some models are only valid if the absolute magnitude of energy loss (as calculated by DELTAE) is larger than a minimal value, such as Moliere multiple scattering.

Default value is 0.025 MeV/n.

EMTRANS : Future switch for photon/electron/positron transportation. Not implemented, default is 0 (off).

EXTSPEC : Specifies if external source file to be read with **USECBEAM** is an array of monoenergetic beams (**EXTSPEC** 0) or a binned spectrum (**EXTSPEC** 1). If a binned spectrum, then the energies are assigned to the lower energy of that bin. The last bin marks the endpoint of the energy spectrum, and its weighting value is thus ignored.

This card is optional, and default is 0 (no spectrum). The card has only effect if **USECBEAM** is called *after* **EXTSPEC**.

HIPROJ : If **JPART0**=25 then the of this card 2 arguments specifies the number of nucleons A and the charge Z of the beam particles. Z has to be larger than 2, since nucleons with $Z = 1$ and $Z = 2$ have their own particle ID **JPART0**, as listed in table 7.2.

If this card is omitted, then default values are $Z = 6$ and $A = 12$.

JPART0 : Optional card. 1 argument that specifies the primary particle ID, see table 7.2. If **JPART0**=25, then the **HIPROJ** card can be called as well. If **HIPROJ** is not called, then carbon-12 is assumed by default.

If **JPART0** is not specified, then the default particle 2 (protons) are assumed.

MAKELN : Optional card. 1 argument that invokes phase-space output of all secondary neutrons with energies below 14.5 MeV which are created within the target: 1 output is written to *for028*, see section 4.1, 0 no output of secondary neutrons.

Default value is 0.

MSCAT : Optional card. 1 argument that switches between the available types of multiple scattering: 1 for Gaussian- and 2 for Moliere-type multiple scattering.

Moliere multiple scattering (2) is activated by default.

NEUTRLCUT : Optional card. Lower energy cut-off value for neutron transport in MeV. Must be lower than 14.5 MeV but larger or equal 0.0 MeV. (This was previously called **OLN** in older versions of **SHIELD-HIT**.)

Default value is 0.0. In this case, neutrons are transported until they are absorbed or exit the simulation universe.

NSTAT : Optional card. 2 arguments that specify the requested total number of primary particles to be simulated and the number of transported primaries after which an intermediate save to the scorers will be invoked, respectively.

Default values are 10000 primary particles simulated, and intermediate saving after 5000 transported primary particles.

NUCRE : Optional card. 1 argument that switches nuclear reactions on or off: 1 with nuclear reactions; 0 all nuclear reactions are turned off. (This card was called **INUCRE** in earlier versions of SHIELD-HIT.)

Default value is 1, i.e. nuclear reactions are activated.

RIFIMC : Optional card. 1 argument that specifies the mode of the ripple filter (RIFI) simulation: 0 for (most realistic) modulus RIFI version retaining the spatial period information of the RIFI; 1 for Monte Carlo sampling of RIFI material thickness—useful when generating SPC files (see section 3.5.6).

Default value is 0.

RIFITRANS : Optional card. 1 argument that specifies the interpretation of data in the Input files loaded with the **USERIFI** card: 0 lists of material thicknesses; 1 lists of vacuum thicknesses.

Default value is 0.

RNDSEED : Optional card. 1 argument that specifies the random number seed. Using the same seed for an identical simulation should yield the same results.

Default value is 89736501.

STRAGG : Optional card. 1 argument that switches between the available models for energy straggling: 1 for Gaussian- and 2 for Vavilov-type energy straggling; 0 for no straggling.

Vavilov straggling (2) is activated by default.

TMAX0 : Optional card. 2 arguments that specify the initial energy of the primary particle in MeV/nucleon and the energy spread in MeV/nucleon, expressed as one standard deviation of a Gaussian distribution (1σ).

Default values are 250 MeV/nucleon and zero energy spread.

USECBEAM : Optional card. 1 argument that specifies the filename of an optional external beam source file. See section 3.3.2.

No file name is specified by default

USEPARLEV : Optional card. 1 argument that specifies the file name of an external file. The file may contain model parameter values that override the default **PARLEV** model parameters. See section 3.3.3.

No file name is specified by default

USERIFI : Optional card. 1 argument that specifies the file name of an optional external ripple filter file. See section 3.3.4.

No file name is specified by default

WRITEFILE : Optional card. A list of 12 flags that toggles generation of various output files. The first flag toggles transportation of fast neutrons (> 14.5 MeV). A flag is activated with **y** and deactivated with **n**. This is described further in section 4.2.

By default fast neutron transportation is turned on, and no output files are generated, i.e. **ynnnnnnnnnnnn**.

3.3.2 USECBEAM - External source file

If a file name is specified, e.g. *sobp.dat*, a projectile will be sampled from the file. Beam settings specified in the input file *beam.dat* such as BEAMSIGMA and TMAX0 will be overridden. Example of a typical beam file, which was generated from a hadron therapy carbon ion control file:

```
*ENERGY(GEV) X(CM) Y(CM) FWHM(cm) WEIGHT
0.270550 -1.00 1.00 0.48 9.3700e+06
0.270550 -0.90 1.00 0.48 9.3700e+06
0.270550 -0.80 1.00 0.48 9.3700e+06
0.270550 -0.70 1.00 0.48 9.3700e+06
0.270550 -0.60 1.00 0.48 9.3700e+06
0.270550 -0.50 1.00 0.48 9.3700e+06
0.270550 -0.40 1.00 0.48 9.3700e+06
0.270550 -0.30 1.00 0.48 9.3700e+06
0.270550 -0.20 1.00 0.48 9.3700e+06
0.270550 -0.10 1.00 0.48 9.3700e+06
(...)
```

Notice that the energy is in GeV to maintain compatibility with other Monte Carlo codes. Lines starting with * are ignored and can be used for comments.

If 6 columns are specified, following format is assumed:

```
*ENERGY(GEV) X(CM) Y(CM) FWHMx(cm) FWHMy(cm) WEIGHT
0.270550 -1.00 1.00 0.48 0.44 9.3700e+06
0.270550 -0.90 1.00 0.48 0.44 9.3700e+06
0.270550 -0.80 1.00 0.48 0.44 9.3700e+06
0.270550 -0.70 1.00 0.48 0.44 9.3700e+06
0.270550 -0.60 1.00 0.48 0.44 9.3700e+06
0.270550 -0.50 1.00 0.48 0.44 9.3700e+06
0.270550 -0.40 1.00 0.48 0.44 9.3700e+06
0.270550 -0.30 1.00 0.48 0.44 9.3700e+06
0.270550 -0.20 1.00 0.48 0.44 9.3700e+06
0.270550 -0.10 1.00 0.48 0.44 9.3700e+06
(...)
```

The FWHM parameter is the Gaussian FWHM (and not sigma), contrary to the BEAM-SIGMA card¹. However, if FWHM is negative (both x and y in the 2-d case), then a box with similar side length is generated. If only one of the two values are negative in the 2-d case, then a flat circular distribution is assumed, with FWHMy as the circular radius. FWHMx is currently ignored, but may be used for ellipsoid beamlets in future.

3.3.3 USEPARLEV - External parameter file

The predefined values of the PARLEV parameters can be redefined by those given in the optional PARLEV input file thereby overwriting the carefully selected and benchmarked default values. Further information on PARLEV parameters can be found in 7.7 and the default values are listed in Table 7.6.

¹Thereby compability to FLUKA can be maintained.

It is recommended to use the default values. They have been tested against available data. Most likely you want to use the default values to remove the `USEPARLEV` card in the *beam.dat* file entirely.

Example of typical `PARLEV` input file with two columns. The first column specifies the ID of a `PARLEV` parameter while the second specifies its new value (either space or tab separated).

```
* ID Value
33 1.0
34 1.0
39 1.20
40 1.20
```

Lines starting with `*` are ignored and can be used for comments.

3.3.4 USERIFI - Ripple filter

A ripple filter [29] (RIFI) can be inserted into the beam at an arbitrary position. Shown below are the first lines of a typical (1D) ripple filter file:

```
* The start position of ripple filter in Z direction
* must coincide with a geometrical area.
* This number is in units of cm.
-1.30
*Y coordinate    Z displacement    both in mm (!)
-0.797653867871    2.99843308865
-0.797583286279    2.96589497459
-0.790976849238    2.92032749859
-0.790878035008    2.8747741389
-0.784271597967    2.8292066629
-0.771129305477    2.77060982496
-0.771072840203    2.74457933371
-0.764466403162    2.69901185771
-0.751366459627    2.6599378882
-0.744760022586    2.6143704122
-0.738167701863    2.57531055901
```

SHIELD-HIT12A assumes by default that the RIFI file lists the thickness of the ripple filter as a function of position. If a displacement in vacuum is stated then the `RIFITRANS` card must be set in *beam.dat*.

Line breaks and lines starting with `*` are ignored and can be used for comments. The ripple filter can either be one or two dimensional. If the ripple filter is one dimensional, the periodic structure changes along the Y axis only. The structure is repeated over the entire ripple filter by applying the modulo operator on the y-position of the incoming particle. Alternatively, the structure can be sampled by a random number using the `RIFIMC` card in *beam.dat*.

The data must be either space or tab delimited and the maximum number of steps is set to 200. The size of the ripple filter is specified in the *geo.dat* file. In the example shown below, the RIFI is represented by the body RPP 3. The side facing closest to the beam **must match** the starting position in the aforementioned ripple filter file to activate the periodic structure. Successful application of the ripple filter will be stated at run time when the first particle hits

the ripple filter surface.

Example: `geo.dat` with ripple filter (body and zone number 3).

```

0      0      C-12 on water with RiFi  21.10.2011
RPP    1    -100.0    100.0    -100.0    100.0    0.0    100.0
RPP    2    -150.0    150.0    -150.0    150.0    -150.0    150.0
RPP    3     -8.0     8.0     -8.0     8.0    -1.300    -1.0
RPP    4    -200.0    200.0    -200.0    200.0    -200.0    200.0
END
TAR    1      +1
AIR    2      +2     -1     -3
RIF    3      +3
OUT    4      +4     -2
END
1      2      3      4
1      2      3      0

```

Additional information about the implementation of the ripple filter can be found in [30] and [10].

A 2-dimensional ripple filter can be specified as well. SHIELD-HIT12A detects this automatically if the specified file contains three columns of data instead of two. The supplied file should be a mesh of data, describing a **half period** of the ripple filter. Either the X coordinate can increment fastest followed by the Y position, or vice versa, yet the data must be ordered in increasing order. The third column gives the thickness in Z direction. All units are in mm, except for the value which designates the start position of the ripple filter, which is in cm, just as in the 1-dimensional case. Line breaks and lines starting with * are ignored. The maximum number of allowed lines is hard coded to be 200x200 steps (i.e. 40.000 data points). The table below shows an excerpt of a 2D ripple filter file example:

```

* 2-D ripple filter example.
* The start position of ripple filter along Z axis
* must coincide with a geometrical area.
* This number is in units of cm.
-1.30
*X coordinate  Y coordinate  Z displacement    all in mm (!)
0 0 3.6
0.01 0 3.6
0.02 0 3.6
0.03 0 3.6
0.04 0 3.6
0.05 0 3.6
0.06 0 3.6
(...many lines skipped here...)
0.73 0.01 1
0.74 0.01 0.95

0 0.02 3.6
0.01 0.02 3.6

```

```
0.02 0.02 3.6
0.03 0.02 3.6
(...many lines skipped here...)
0.69 0.74 0
0.7 0.74 0
0.71 0.74 0
0.72 0.74 0
0.73 0.74 0
0.74 0.74 0
```

3.4 *geo.dat* - Geometry

The SHIELD-HIT code uses known Combinatorial Geometry (CG), similar to the Monte Carlo particle transport codes MORSE, FLUKA.

The geometry is defined by a number of so called *bodies* which are primitive objects such as boxes and cylinders, which by Boolean logic are combined to *zones*. For instance a square slab with a round hole can be constructed using two primitives, i.e. a box subtracted with a cylinder. Similarly a hollow box can be described by a larger box subtracted with a smaller box inside it.

The zones must then be assigned to a material identifier, which is an integer defined by the user in the material definition file *mat.dat*, see section 3.2. Every point in the simulated universe must be assigned to a zone, else errors may occur. So, if a hollow box is constructed, one must remember also to assign the void inside the box to a material code (e.g. internal vacuum 1000). One should also understand that a single body cannot directly be assigned to a material, but must be assigned to a zone first, even if the zone just consists of a single body.

The definition of the bodies, zones and the material assignments are set in the *geo.dat* file, in exactly that order.

Anywhere within *geo.dat* lines starting with *** are skipped.

3.4.1 Title

The first line of *geo.dat* contains two integers JDBG1, JDBG2 and a string of characters describing the title of the geometry. The format of the title card is:

- 5 columns (1-5): integer JDBG1, see below
- 5 columns (6-10): integer JDBG2, see below
- 10 columns (11-20): unused
- 60 columns (21-80): any title of the geometry

The first integer JDBG1 selects whether the file *for017* containing the geometry debugging information should be kept (0) or deleted (1) after the geometry parser was initialized. The second integer JDBG2 describes the lower cutoff value of transportation step size in powers of 10, i.e. $10^{-|\text{JDBG2}|}$. If set to zero, the minimum step size is set to 10^{-4} cm. If any of the numbers are omitted, they are interpreted as (0). The fields for both numbers are five columns wide.

3.4.2 Bodies

A list of bodies describing the geometry follows the title card explained in the section before. A body is a primitive object with a given geometrical configuration. The available primitive objects are listed section 7.5. A “top level” body (which is mandatory in the MC code FLUKA) can optionally be added which contains all other bodies, which will be the simulation universe, but this is not strictly necessary in SHIELD-HIT12A. There are no requirements to how this body must be named or where it is placed in the body list, but for convenience it can be either the first or last body. The list of bodies must be terminated with the **END** card. The format of a body card is:

- 2 columns (1-2): unused, e.g. two empty spaces
- 3 columns (3-5): name of the body (see sec. 7.5)
- 1 column (6): unused

- 4 columns (7-10): an integer describing the number of the body
- 6 x 10 columns (11-70): 6 arguments describing the body as mentioned in 7.5.

A continuation card may be necessary, depending on what body is described. A continuation card must be blank in the first 10 columns (i.e. no name and no body number). The body section must be terminated with an **END** card.

3.4.3 Zones

The zone description follows afterwards. In order to construct different zones out of the bodies defined at the beginning of *geo.dat*. Boolean logic is used. Examples of how to construct zones out of two bodies are shown in figure 3.2.

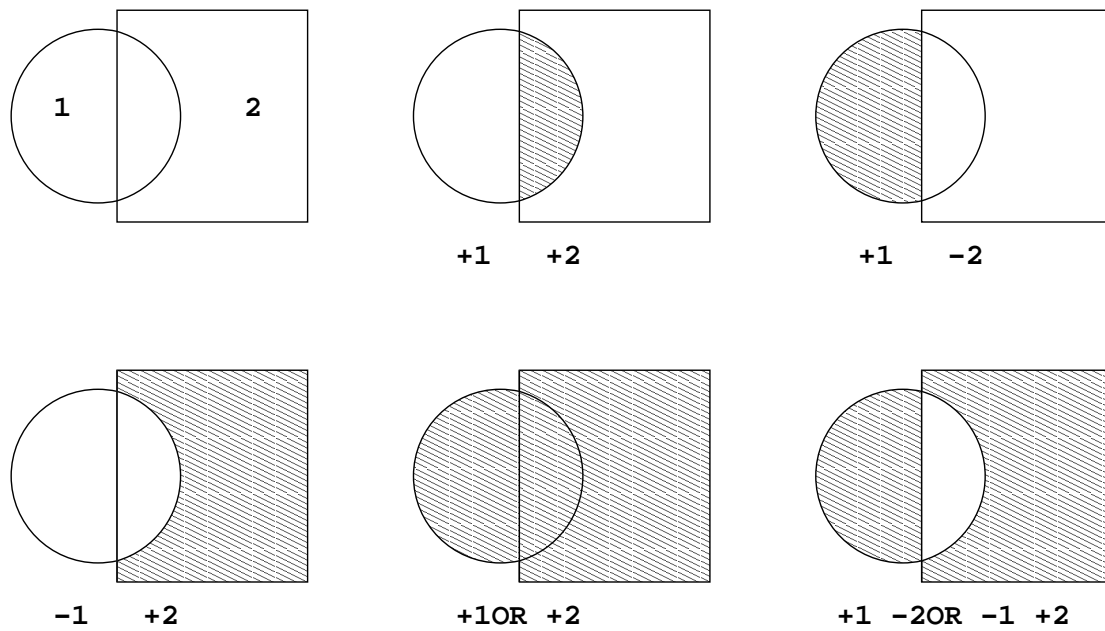


Figure 3.2: Two bodies (marked as 1 and 2) describing 5 different zones.

A line in the zone description consists of:

- 2 columns (1-2): unused, e.g. two empty spaces
- 3 columns (3-5): name of the zone
- 5 columns (6-10): an optional integer describing the number of the zone
- 9 x (2+5) columns (11-72): where each set of 9 columns are composed of two components. The two first columns hold the Boolean operator **OR** or are left blank if needed. The following 5 columns contain a signed (+,-) integer body number. It is essential that the alignment is strictly kept.

If more than 9 bodies are used for the zone description additional continuation cards can be added directly afterwards. These continuation lines must not have a zone name, i.e. there must be no entry at columns 3 - 5. The zone section must also be terminated with **END** card.

Example:

```

* <-><---> <--->OR<--->OR<--->OR<--->OR<--->OR<--->OR<--->OR<--->OR<--->
001      +1      -2      -3      -4      -5
002      +2
003      +3
004      +4
005      +5
006      +6      -1
END
\begin{Verbatim}

```

It is important that any point (except for zone boundaries) where a particle may go, is described by exactly one zone, no more, no less.

\SHA\ may produce unexpected or random behavior if particles cross areas where two zones overlap, or a zone which is not defined. To some extent \SHA\ will warn the user in such cases, but \SHA\ may not catch all cases, therefore a careful assignment of zones is essential.

\subsection{Media}

Finally, what follows is the assignment of media to the various zones. External vacuum behaves like a black hole, any particle hitting it will be disintegrated and not transported further. The ‘‘top level’’ zone must be assigned to external vacuum (id = 0).

Inside this zone you can place a zone with internal (i.e. real physical) vacuum (id = 1000) or any other material. In internal vacuum regular particle transport takes place, as opposed to the external vacuum. Even if not strictly mandatory, the external vacuum should encompass the entire system is to prevent infinite loops where stray particles are transported forever, and to avoid unintended behavior of the code. The media assignment is done in two equally formatted lists. The first list is a list of zones:

```

\begin{itemize}
  \item 14x5 columns (1-70): list of integer zone numbers
\end{itemize}

```

The second list are the material numbers:

```

\begin{itemize}
  \item 14x5 columns (1-70): list of integer media identifiers
\end{itemize}

```

If one line is not enough, then it is simply continued on the following line, until all zones are described. All zones must be assigned to medium, thus, both lists must be equally long.

A simple example of the input geometric file \file{geo.dat} is given below.\

```

\begin{Verbatim}[frame=single, label=\fbox{\textbf{Example: water tank surrounded by air}}

```

```

0      0      simple water tank  20.09.2011
RPP    1    -100.0    100.0    -100.0    100.0      0.0    100.0
RPP    2    -150.0    150.0    -150.0    150.0    -150.0    150.0

```



```

RPP      3      -200.0      200.0      -200.0      200.0      -200.0      200.0
END
TAR      1      +1
AIR      2      +2      -1
OUT      3      +3      -2
END
      1      2      3
      1      2      0

```

A more complex example is a model of the beam line at the cyclotron facility in Clatterbridge, UK. It features two zones with external vacuum, where the second one is used to collimate the beam:

```

      0      0      COO beam line Clatterbridge 13.05.2011
RCC      1      0.0      0.0      -25.6      0.0      0.0      1.00
      15.0
RCC      2      0.0      0.0      -25.6      0.0      0.0      1.00
      0.3
RCC      3      0.0      0.0      -22.6      0.0      0.0      0.003
      15.0
RCC      4      0.0      0.0      -0.66      0.0      0.0      0.660
      0.286
RCC      5      0.0      0.0      0.0      0.0      0.0      0.002
      15.0
RCC      6      0.0      0.0      5.0      0.0      0.0      0.005
      15.0
RCC      7      0.0      0.0      -30.0      0.0      0.0      35.005
      15.0
RCC      8      0.0      0.0      27.0      0.0      0.0      0.001
      20.0
RCC      9      0.0      0.0      55.0      0.0      0.0      1.00
      20.0
RCC     10      0.0      0.0      55.0      0.0      0.0      1.00
      2.0
RCC     11      0.0      0.0      115.00      0.0      0.0      0.002
      20.0
RCC     12      0.0      0.0      115.002      0.0      0.0      0.002
      20.0
RCC     13      0.0      0.0      175.0      0.0      0.0      7.80
      20.0
RCC     14      0.0      0.0      175.0      0.0      0.0      7.80
      1.70
RCC     15      0.0      0.0      182.80      0.0      0.0      0.80
      20.0
RCC     16      0.0      0.0      182.80      0.0      0.0      0.80
      1.5
RCC     17      0.0      0.0      5.005      0.0      0.0      178.995
      25.0
RCC     18      0.0      0.0      184.0      0.0      0.0      30.0
      20.0

```

```
RCC  19      0.0      0.0    -100.0      0.0      0.0      500.0
      100.0
RCC  20      0.0      0.0    -200.0      0.0      0.0     1000.0
      200.0
RCC  21      0.0      0.0     -26.0      0.0      0.0        1.00
      15.0
RCC  22      0.0      0.0     -26.0      0.0      0.0        1.00
      0.3
END
TAR   1    +18
BR1   2     +1    -2
BR2   3     +9   -10
BR3   4    +13   -14
BR4   5    +15   -16
HL1   6     +2
HL2   7    +10
HL3   8    +14
HL4   9    +16
SF1  10     +3
SF2  11     +5
STP  12     +4
KWD  13     +6
VAC  14     +7    -1    -3    -4    -5    -6   -21
WHL  15     +8
ICM  16    +11
ICA  17    +12
CAR  18    +17    -8    -9   -11   -12   -13   -15
OAR  19    +19    -7   -17   -18
OUT  20    +20   -19
IBH  21    +21   -22
IHL  22    +22
END
  1   2   3   4   5   6   7   8   9  10  11  12  13  14
15  16  17  18  19  20  21  22
  1   2   2   2   2   2   2   3   3   9   9   2   5 1000
  8   6   4   3   3   0   0 1000
```

%% TODO: Nikolai: two outer vaccuum (21 and 22) is this correct?

3.5 *detect.dat* - Auxiliary scoring

Different geometrical scoring types, called *estimators*, are available for auxiliary scoring: ZONE, DZONE, CYL, MSH, PLANE DCYL, DMSH, DMSHZ and DPLANE. Each estimator can be invoked with a scoring card and possible following cards in the file *detect.dat*. Most estimators need additional specification of a *detector*, which controls what kind of quantity is scored. The available detectors are described afterwards in section 3.5.7.

The output of each estimator is written, in binary form, to a file, which is specified by the user. The post-processing script *shield.detect2ascii*, described in section 5.2.1, translates the binary output (of some estimators) to human readable ASCII format.

3.5.1 ZONE - Scoring by zone

ZONE : Scoring by zones w_i . Zones are defined from bodies in *geo.dat*. A single or a range of zones can be specified. Arguments are:

1. w_i : first zone of a range of zones to be scored.
2. w_j : optional last zone of a range of zones. If this is not specified, then only a single zone w_i will be scored. Else every zone within the interval $[w_i, w_j]$ will be scored.
3. unused.
4. **JPART**: particle type. See list in section 7.2. Setting this value to -1 scores all particles. If heavy ions are scored (**JPART**=25) then a continuation card is required, where the first argument is an integer representing the particle charge Z , and the second argument is an integer representing the particle mass A .
5. **DETECTOR**: quantity that should be scored. A list of available detectors is given in section 3.5.7.
6. *outputfile*: output file name.

Analogue to CYL and MSH, a continuation card is only required if **JPART**=25 or if **DETECTOR** is specified as LETFLU (and only LETFLU). Arguments for the continuation card are:

1. Z : an integer specifying the charge of the particle to be scored.
2. A : an integer specifying the number of nucleons of the particle to be scored.
3. **MEDIUM**: an integer specifying the medium for LET weighted scoring. If left unspecified, the default is the medium material that the bin is made of. Any other integer from the medium list specified in *mat.dat* is valid here. In that case, scoring will apply the stopping powers of the medium specified here. Transport will irrespective of the scoring always be done in the as medium specified in *geo.dat*. This is useful for calculating stopping power ratios needed by various cavity theories in radiation dosimetry.
4. unused.
5. unused.
6. unused.

3.5.2 CYL - Cylindrical scoring

CYL : Cylindrical scoring. Cylindrical coordinates are used. Currently, scoring can only be done along the Z axis. This card requires a second succeeding card. Arguments are:

1. R_{min} : inner radius of scoring cylinder. If $R_{min} > 0$ then the scoring volume is a cylinder shell.
2. θ_{min} : minimum value of angular segment.
3. Z_{min} : start position of cylindrical scoring in Z direction.
4. R_{max} : outer radius of scoring cylinder.
5. θ_{max} : maximum angle of angular segment. If this value is set to number larger than 2π then θ_{max} will be fixed at 2π , covering the entire disc.
6. Z_{max} : end position of cylindrical scoring in Z direction.

Arguments for the second card:

1. R_{bin} : number of bins in radial direction.
2. θ_{bin} : number of angular bin segments.
3. Z_{bin} : number of bins in Z direction.
4. **JPART**: particle type. See list in section 7.2. Setting this value to -1 scores all particles. If heavy ions are scored (**JPART**=25) then a third continuation card is required, where the first argument is an integer representing the particle charge Z , and the second argument is an integer representing the particle mass A .
5. **DETECTOR**: quantity that should be scored. A list of available detectors is given in section 3.5.7.
6. *outputfile*: output file name.

Analogue to **ZONE** and **MSH** a continuation card is only required if **JPART**=25 or if **DETECTOR** is specified as **LETFLU** (and only **LETFLU**). Arguments for the third card are:

1. Z : an integer specifying the charge of the particle to be scored.
2. A : an integer specifying the number of nucleons of the particle to be scored.
3. **MEDIUM**: an integer specifying the medium for LET weighted scoring. If left unspecified, the default is the medium material that the bin is made of. Any other integer from the medium list specified in *mat.dat* is valid here.
4. unused.
5. unused.
6. unused.

3.5.3 MSH - Cartesian scoring

MSH : Cartesian mesh scoring card. This card requires a second succeeding card. Arguments for the first card:

1. X_{min} : lowest X position.
2. Y_{min} : lowest Y position.
3. Z_{min} : lowest Z position.
4. X_{max} : highest X position.
5. Y_{max} : highest Y position.
6. Z_{max} : highest Z position.

Arguments for the second card:

1. X_{bin} : number of bins in X direction.
2. Y_{bin} : number of bins in Y direction.
3. Z_{bin} : number of bins in Z direction.
4. JPART: particle type. See list in section 7.2. Setting this value to -1 scores all particles. If heavy ions are scored (JPART=25) then a third continuation card is required, where the first argument is an integer representing the particle charge Z , and the second argument is an integer representing the particle mass A .
5. DETECTOR: quantity that should be scored. A list of available detectors is given in section 3.5.7.
6. *outputfile*: output file name.

Analogue to ZONE and CYL a continuation card is only required if JPART=25 or if DETECTOR is specified as LETFLU (and only LETFLU). Arguments for the third card are:

1. Z : an integer specifying the charge of the particle to be scored.
2. A : an integer specifying the number of nucleons of the particle to be scored.
3. MEDIUM: an integer specifying the medium for LET weighted scoring. If left unspecified, the default is the medium material which the bin is made of. Any other integer from the medium list specified in *mat.dat* is valid here.
4. unused.
5. unused.
6. unused.

3.5.4 PLANE - Scoring by 2D plane

PLANE : Scoring by a single 2-D infinite plane s , independently of *geo.dat*. The plane is specified by a random point which must lie in the plane, and the normal vector of that plane. The normal can either be specified in this card, or - if omitted - the normal vector will automatically be calculated from the beam direction specified by the **BEAMTHETA** and **BEAMPHI** cards in *beam.dat*.

Arguments are:

1. S_x : X coordinate of point in plane.
2. S_y : Y coordinate of point in plane.
3. S_z : Z coordinate of point in plane.
4. n_x : x component of normal vector.
5. n_y : y component of normal vector.
6. n_z : z component of normal vector.

Arguments for the second card:

1. unused.
2. unused.
3. unused.
4. **JPART**: particle type. See list in section 7.2. Setting this value to -1 scores all particles. If heavy ions are scored (**JPART**=25) then a third continuation card is required, where the first argument is an integer representing the particle charge Z , and the second argument is an integer representing the particle mass A .
5. **DETECTOR**: quantity that should be scored. A list of available detectors is given in section 3.5.7.
6. *outputfile*: output file name.

Analogue to **CYL** and **MSH**, a continuation card is only required if **JPART**=25 or if **DETECTOR** is specified as **LETFLU** (and only **LETFLU**). Arguments for the continuation card are:

1. Z : an integer specifying the charge of the particle to be scored.
2. A : an integer specifying the number of nucleons of the particle to be scored.
3. **MEDIUM**: an integer specifying the medium for LET weighted scoring. If left unspecified, the default is the medium material that the bin is made of. Any other integer from the medium list specified in *mat.dat* is valid here. In that case, scoring will apply the stopping powers of the medium specified here. Transport will irrespective of the scoring always be done in the as medium specified in *geo.dat*. This is useful for calculating stopping power ratios needed by various cavity theories in radiation dosimetry.
4. unused.
5. unused.
6. unused.

3.5.5 DZONE, DCYL, DMSH and DPLANE - Differential scoring

DZONE, DCYL, DMSH and DPLANE : Differential zone, cylindrical and mesh scoring. Same as ZONE, CYL and MSH, respectively, but with an extra card that specifies the differential dimension of the scoring. The differential scoring card must be inserted before the heavy ion selection card, if that is used as well.

Arguments for the extra card:

1. Δ_{start} : Start value of differential binning.
2. Δ_{stop} : Stop value of differential binning.
3. Δ_n : Number of bins. If this value is negative, then binning is done in logarithmic steps, and the Δ_{start} parameter must be larger than zero.
4. internal use, keep this field clear.
5. **Type**: Type of differential scoring. Currently, ENERGY, LET and ANGLE are available, being differential in MeV/n, MeV/cm or radians, respectively.
6. unused.

The optional third card for CYL and MSH is the fourth card when using DCYL and DMSH.

If applying differential scoring on pions, their energies will be converted to specific energies (MeV/n) by multiplying with 0.1498 (π^+ and π^- mass divided by proton mass)².

²In the case of π^0 it is 0.1448 which the π^0 mass divided by proton mass. However, SHIELD-HIT12A does not transport π^0 as it decays instantaneously.

3.5.6 DMSHZ and DCYLZ - Spectral files in TRiP98 format

DMSHZ and DCYLZ : Creates a `.spc` file for TRiP98 [31] format using either mesh or cylindrical scoring. The SPC format is described in

<http://bio.gsi.de/DOCS/TRiP98/PRO/DOCS/trip98fmtspc.html>.

DMSHZ and DCYLZ is similar to DMSH DCYL, but requires the special detector SPC in the DETECTOR field of the second card. The mandatory third card specifies differential dimensions of the scoring:

1. Δ_{start} : Start of differential binning in energy (MeV/A).
2. Δ_{stop} : Stop of differential binning in energy (MeV/A).
3. Δ_n : Number of bins. If this value is negative, then binning is done in logarithmic steps, and the Δ_{start} parameter must be larger than zero.
4. Z_{peak} : Peak position for placement of scoring grid along Z. If negative the peak position will be estimated by SHIELD-HIT12A in CSDA approximation, assuming material NUMMED=1.
5. Type: ENERGY, only.
6. unused.

For DMSHZ and DCYLZ scoring, the first medium NUMMED=1 specified in *mat.dat* can be any material, but most likely one may use water (e.g. for generating proper TRiP spc files). In addition, the SPC writer must be the last detector specified in *detect.dat*, otherwise SHIELD-HIT12A may behave in an undefined way.

The peak position Z_{peak} is needed to place the adaptive scoring grid in Z direction which has the highest resolution at the Bragg-peak, and will be more coarse the further away the scoring is from the peak position. If a negative number is given, SHIELD-HIT12A will estimate the peak position using the continuous slowing down approximation (CSDA) in water. This may be insufficient when a true beam line is modelled, which might significantly shift the true Z_{peak} upstream. In this case, the center of the grid can be placed manually by specifying a positive Z_{peak} .

The Z_{max} position specified in the first card of the DMSHZ estimator will be used by SHIELD-HIT12A for setting up the Z interval wherein the Bragg-peak position will be found. For the scoring of the spectrum, however, Z_{max} will be then overwritten by SHIELD-HIT12A by twice the Z_{peak} position. Currently only $Z_{min} = 0$ is possible.

The specified number of Z bins, Δ_n , is not propagated to the output SPC-files, but internally used by SHIELD-HIT12A to determine the Bragg peak position along the Z axis. This should not be confused with Z_{peak} specified by the user, as the result of Bragg-peak position calculation is only available after transportation. The exact position of the Bragg-peak location is then written in the SPC file, needed by TRiP.

When handling divergent beams, the distance to the isocenter becomes important. For DDD and SPC file generation, a constant source to surface distance (SSD) is assumed, thus the isocenter is at a fixed position in the target, irrespectly of the penetration depth of the ions. However, in treatment planning in TRiP the isocenter is located in the treatment volume, which means varying SSDs. While unimportant for non-diverging beams, the solution for divergent beams can be twofold.

1. Newer versions of TRiP provide a new parameter for the `scancap` command which solves this issue.

2. Older versions of TRiP one should recalculate a new set of base data for each field, which uses a proper SSD and isocenter depth corresponding to this particular field.

Contrary to the behaviour of DMSH and DCYL, DMSHZ and DCYLZ will score all (primary or secondary) pions and antiprotons as protons with the same energy per mass. The reason is that TRiP cannot handle pion or antiprotons per default. Yet one can approximate that the antiproton (when treating the secondaries from annihilation separately) resembles a proton in terms of dose deposition and stopping power. Pions have also unity charge, but are 7 times lighter than protons. A 70 MeV pion is therefore scored as a 10 MeV proton, since these two particles have comparable stopping powers at these energies.

3.5.7 Available detectors for scoring

Different detector types are available for the auxiliary scoring. For each geometrical estimator invoked in the file `detect.dat` one type of detector has to be specified for **DETECTOR** in the second card. But not every geometrical estimator type can be combined with each of the available detectors. A detector type may itself invoke an additional card. The output is given in SHIELD-HIT12A standard units (see table 7.1). The quantities that scale with particle number are normalized per simulated primary particle. The available detectors are:

ENERGY : Absolute energy deposited in each bin.

FLUENCE : Track length fluence in each bin ϕ .

CROSSFLU : Absolute number of particles entering each bin.

DOSE : Energy deposited in each bin divided by total mass (volume \times density) of the bin.

LETFLU : Sum of the products between fluence and unrestricted linear energy transfer (LET) over particle type i ,

$$\bar{\phi}_{\text{LET}} = \sum_i \phi_i(E, Z) \frac{dE}{dx}(E, Z)_i, \quad (3.1)$$

for a medium specified by card 3, argument 3 which might be different from the material of the bin. This argument is a number m that corresponds to the m^{th} material in `mat.dat`. The product for one specific particle i can be selected with arguments 1 and 2 in card 3. All particles are considered if these two arguments are left blank.

DLET : Dose (D) averaged unrestricted stopping power (LET) in MeV/cm. For each particle i :

$$\overline{LET}_D = \frac{\sum_i D_i(E, Z) \frac{dE}{dx}(E, Z)_i}{\sum_i D_i(E, Z)}. \quad (3.2)$$

TLET : Track length (τ) - averaged unrestricted stopping power (LET) in MeV/cm. For each particle i :

$$\overline{LET}_\tau = \frac{\sum_i \tau_i(E, Z) \frac{dE}{dx}(E, Z)_i}{\sum_i \tau_i(E, Z)}. \quad (3.3)$$

which equals the fluence (ϕ) averaged unrestricted stopping power, since $\phi = \tau/V$.

AVG-ENERGY : Track length averaged energy. For each particle i :

$$\bar{E}_\tau = \frac{\sum_i E_i(Z) \tau_i(E, Z)}{\sum_i \tau_i(E, Z)}. \quad (3.4)$$

AVG-BETA : Track length averaged relativistic velocity $\beta = v/c$. For each particle i :

$$\bar{\beta}_\tau = \frac{\sum_i \beta_i(Z) \tau_i(E, Z)}{\sum_i \tau_i(E, Z)}. \quad (3.5)$$

SPC : Only available for DMSHZ. See section 3.5.6.

ALANINE : Response-equivalent dose for the alanine detector. The medium in which this detector is applied must be alanine (ICRU ID = 105) in order to make sense. For each charged particle i :

$$D_{\text{al,photon}} = \sum_i \eta(Z_i, E_i) D_{i,\text{alanine,ion}}. \quad (3.6)$$

The relative effectiveness η is calculated using an implementation [32] of the Hansen and Olsen model [33,34] which only takes pions, kaons, and ions with $Z < 7$ into account. For heavier ions ($Z > 6$) η is set to 1.0. For energies below 1 keV/nucleon η is fixed to 0.3 for all ions.

COUNTER : A counter which counts the number of particles crossing a **PLANE** estimator.

3.5.8 Scoring examples

This example shows how to score a 1-D energy deposition of all particles in a cylindrical volume which is 20 cm long (from 35 cm to 55 cm). The output will be written into the file *foo*. In addition, a 2-D fluence map ($1 \times 20 \times 300$ bins) of carbon-12 ions ($Z = 6$ and $A = 12$) is generated and written to the file *bar*. Finally, the average energy deposited by all particles per primary particle in the zones 2 to 12 defined in *geo.dat* is outputted to *foobar*.

```
*---0--->---1--->---2--->---3--->---4--->---5--->---6--->
CYL          0.0      0.0      35.0      20.0      7.0      55.0
              1        1        300        -1      ENERGY      foo
MSH          0.0      0.0      0.0      10.0      10.0      10.0
              1        20       300        25      FLUENCE      bar
              6        12
ZONE         2        12              -1      ENERGY      foobar
```

The second example shows how to produce two .spc files in **TRIP98** [31] format. The first file *lin.spc* contains the particle spectrum at different depths resolved by particle on a linear energy scale while for the second file, *log.spc*, a logarithmic energy scale is used.

```
*---0--->---1--->---2--->---3--->---4--->---5--->---6--->
DMSHZ       -100.0   -100.0      0.0     100.0     100.0     50.0
              1        1      5000        -1        SPC      lin.spc
              0.0     800.0      800        -1      ENERGY
DMSHZ       -100.0   -100.0      0.0     100.0     100.0     50.0
              1        1      5000        -1        SPC      log.spc
              0.1    1000.0     -100        -1      ENERGY
```

Here is an example, how to record the angular histogram of particles crossing a plane. A X-Y plane is placed at $Z = 10.0$ cm. Differential angle scoring is invoked, scoring all from 0.0 to 5.0 mrad in 20 bins.

```
*---0--->---1--->---2--->---3--->---4--->---5--->---6--->
DPLANE       0.0      0.0      10.0      0.0      0.0      1.0
              0.0     0.005      20              -1      COUNTER NB_count4
              ANGLE
```


Chapter 4

Output files

SHIELD-HIT12A provides several output files, depending on the configuration in *geo.dat* and *detect.dat*. Six have auxiliary and information functions and described briefly in section 4.1. These files are:

for017
for024
for025
for026
for027
for028

The other six possible output files are mostly binary objects for further processing using the legacy scoring system of SHIELD-HIT. They are described in more details in the SHIELD-HIT12A developer's guide. A short overview is given in section 4.1. Of these files, only *for033* may be of interest, which is in ASCII format. Programs to convert the binary files to human readable ASCII exist, but are not included in the official distribution, since the new scoring system is encouraged. If you for some reason need it, the conversion scripts are available on request.

The binary files are:

for029
for030
for031
for032
for033
for034

4.1 Auxiliary output files

These files are not intended for further processing. They contain auxiliary information for the user in the manner of man-readable ASCII coding. There are no scripts for handling these files in an parallelized environment.

for017 : File for checking how the input geometric file *geo.dat* was read by the geometric program *gemca.f*. Usually may be deleted.

for024 : Contains input and output data of the current task. Part of these data is duplicated in the output file *for033*, see below. It is recommended to keep this file.

for025 : Diagnostic file. Contains diagnostics, if any, which appears during run of the task.

for026 : Stopping powers, dependencies range-energy, macroscopic cross sections and other useful information for all particles/fragments and all materials in the task. It is recommended to keep this file.

for027 : Multigroup neutron cross sections below 14.5 MeV for all chemical elements in the task.

for028 : May contain secondary neutrons below 14.5 MeV with all individual parameters: the birth point XYZ, direction of fly, kinetic energy. Usually this is empty file. This file can be used as a neutron source for transportation with MCNP [21] like programs.

for029 : Contains track length (TLE) fluences of zones in binary form.

for030 : Contains double differential yields of particles/fragments from the entire surface of a target assembly in binary form.

for031 : Contains energy depositions in target zones in binary form.

for032 : Contains energy depositions in target zones, decomposed according to LET intervals, defined at compilation time. The file is in binary form.

for033 : Human readable ASCII file that includes:

- Total energy depositions in target zones from all particles/fragments plus some additional information.
- Decomposition of the energy depositions according to generations of particles/fragments.
- Distribution of PET isotopes on target zones.

for034 : Contains fluence depositions at the intersection of cylindrical targets. This file is in binary form.

4.2 Management of output with help of special flags

The string of flags `WRITEFILE` in `beam.dat` defines which output files are to be generated, and can toggle fast neutron transport. It may look, as an example, like this: `yyyyyyynnnyyy` (n-no, y=yes)

Separate positions of `writefile` in the input file `beam.dat` mean the following.

```
writefile(1:1) - "n" - neutron transport ( $E_n > 14.5$  MeV) switched off
    "y" - neutron transport ( $E_n > 14.5$  MeV) switched on

writefile(2:2) - not used, reserved

writefile(3:3) - not used, reserved

writefile(4:4) - not used, reserved

writefile(5:5) - not used, reserved

writefile(6:6) - generates the file for033

writefile(7:7) "y" - generates the output file for032 (Flag "n" saves about 40% computer
    time)

writefile(8:8) "y" - controls the estimation of crossing fluence on intersections, valid for
    cylindric target only
    "n" - no the estimation
    "y" - estimation of the ordinary fluence ( $W = 1/|\cos(np)|$ );
    output of the file for034 (cks.bin)
    "z" - estimation of the planar fluence ( $W=1$ );
    output of the file for034 (cks.bin)

writefile(9:9) "y" - printout PET-isotopes to the file for024

writefile(10:10) "y" - generates the file for031

writefile(11:11) "y" - generates the file for030

writefile(12:12) "y" - generates the file for029
```


Chapter 5

Auxiliary programs

5.1 Helper scripts

5.1.1 *shield_clean* - Remove generated files

The *shield_clean* script removes the files *for017 for024 for025 for026 for027 for028 for029 for030 for031 for032 for033*. If a directory is given in the argument, it will change to that directory first.

In Windows distributions, the file is named *shield_clean.bat*.

5.2 Format conversion

5.2.1 *shield_detect2ascii* - Convert binary data to ascii from *detect.dat*

shield_detect2ascii converts the output of an estimator specified in *detect.dat* (see section 3.5). If run in a parallel environment, and multiple outputs exist of the same estimator, then *shield_detect2ascii* will average over all files specified in the argument. *shield_detect2ascii* can only convert output from MSH and CYL estimators.

5.2.2 *shield2fluka* - Convert *geo.dat* to FLUKA format

shield2fluka converts the geometry file to a FLUKA formatted input file. Only the geometry file is converted, not the beam settings or scoring. The produced fluka formatted *output.inp* cannot be run in FLUKA, but it can be read by “SimpleGeo” which is a windows program for viewing geometries in 3D.

See <http://theis.web.cern.ch/theis/simplegeo/>.

5.2.3 *ddd.py* - Convert cylindrical depth dose binning to DDD format

5.2.4 *shield_dEdx* - Prepare an external stopping power table

shield_dEdx generates a properly formatted stopping power table which can be read by SHIELD-HIT12A. *shield_dEdx* requires the *libdedx* software package installed [3,26], which can be downloaded for free at <http://libdedx.sf.net>.

The libdEdx library features access to a range of alternative stopping power tables and a comprehensive list of target materials. An output files is generated according to the nomenclature found in table 7.4. The generated ASCII file is human readable and contains a header with information on creation date, version number etc.

Example of usage:

```
$ ./processing/dEdx/shield_dEdx 1 223  
Preparing stopping power table for PMMA using PSTAR + MSTAR.  
Wrote Lucite.txt.
```

Chapter 6

Parallelization

SHIELD-HIT12A can run in a parallelized environment in several ways, ranging from single PC which features multiple cores, or massive computer clusters. SHIELD-HIT12A runs on CPUs only, currently no GPU code is implemented. In this chapter several ways of parallelizing SHIELD-HIT12A is explained. SHIELD-HIT12A will here run in “embarrassingly parallel” mode, where simple independent jobs are submitted to various cores.

6.1 Torque

Torque is a widely used system for job submission across multiple nodes. Here a very simple tutorial is given for using multiple nodes on a PC running SHIELD-HIT12A processes. It is assumed that Torque and the PBS scheduler is running and correctly configured. Furthermore it is assumed that SHIELD-HIT12A was fully installed.

A simple PBS submission script *rtshield.sh* can look like this:

```
#!/bin/sh
#
#PBS -N SHIELDHIT_JOB
#PBS -M bassler@phys.au.dk
#
shieldhit $1 -N$PBS_ARRAYID
```

Simply create this file in the directory you wish to run. Step into the same directory and run:

```
$ qsub -V -t 0-9 -d . rtshield.sh
```

The `-V` forwards the local environment variables to the submission script. This way we do not need to specify the path of the `shieldhit` command, since it is referred to by the `PATH` environment variable.

The `-t` option specifies the nodes to be run, in this case from 0 to 9, meaning 10 instances in total. The output files will be suffixed accordingly, so they will not overwrite each other.

Finally, the `-d .` option specifies the working directory. `qsub` expects to find all the input files as well as the *rtshield.sh* script here.

You can check the status of your job submission with `qstat`. A good manual on Torque can be found on <http://www.clusterresources.com/torquedocs/index.shtml>.

After the run completes, each output file (which can be merged by the *shield_detect2ascii* script) will be appended with an integer number corresponding to the instance number. Thus in the example above, the output files will be suffixed with integers ranging from 0000 to 0009.

6.2 Condor

Another job submission and scheduling system is CONDOR¹. Here submission can be handled by the *rcshield.py* script.

```
$ rcshield.py -c -M100 examples/simple
```

will run the job found in the *examples/simple* directory 100 times. All results, debugging output is transferred back to the directory, and each result file is suffixed with a number in order to distinguish the results. *rcshield.py* accepts a range of options:

- h** - print all options with a short description.
- f** - should be followed by a comma separated (without space) file list, which will be transmitted along with the usual mandatory files. The usual input files which always will be transmitted are: *mat.dat geo.dat beam.dat Air.txt Water.txt detect.dat*. Additional files could be a ripple filter file *rifi.dat*, source or a files which overrides the default model parameters (see section 3.3.3).
- M** - number of simulations to be performed. Random seed is automatically changed for each run.
- N** - first run will have the number followed by this option. Useful if you want to add more statistics to a previous batch or runs. If the last completed run had number 99, you should set **-N100**.
- a** - after each batch of runs, *rcfluka.py* will wait 30 seconds, for any file transfers to complete. This grace time is necessary if you run any postprocessing scripts or the **-c** option. For small runs and quick testing, the 30 secs may be annoying, and can therefore be deactivated with the **-a** option.
- n** - the jobs are submitted as nice jobs to the cluster, backfilling it if no other tasks are pending. Use this option if you are not in a hurry to get your data.
- c** - clean up some of the condor files which were generated, which are only of interest if something goes wrong. Purpose of this option is to keep your run directory tidy.
- C** - big clean: use this if you want to start from scratch in your directory. Nothing is submitted to the cluster, but the directory is emptied for any data generated by previous runs. DANGEROUS, files which may resemble generated data are deleted as well. Basically anything which matches *_rcshield**, **[0-9][0-9][0-9][0-9]*, *for[0-9][0-9][0-9]* and *fort.24* are removed.
- t** - test script, but do not submit jobs to condor cluster.
- S** - use standard universe instead of vanilla (see condor manual). Standard universe will only work if you have linked in the appropriate condor libraries for checkpointing into the SHIELD-HIT12A executable.

¹<http://research.cs.wisc.edu/condor/>

- o** - followed by a string which will be added to the condor submit file.
- i** - followed by a filename of a file which will be appended to the condor submit script.
- m** - followed by an email address, which will be notified when all jobs are completed (or failed).
- x** - single submission mode. This overrides -M and -m option. Simply one job is submitted, and *rcshield.py* immediately exits after submission. This option is useful if you just want the raw output from a single run, and dont want the *rcshield.py* script block the terminal while running. Entire control is within condor, which will transfer files upon job completion.

Chapter 7

User's reference tables

7.1 Units

Table 7.1 provides an overview of the standard units in SHIELD-HIT12A. They apply at any time unless explicitly stated otherwise.

Table 7.1: Table of default units in SHIELD-HIT12A.

Value	Default unit
Distance; position	cm
Track length density; fluence	cm^{-2}
Volume	cm^3
Density	g/cm^3
Atomic density	atoms/cm^3
Projectile energy (ions)	MeV/nucleon
Projectile energy (non-ions)	MeV
Deposited energy	MeV
Dose	MeV/g
Stopping power; LET	MeV/cm
Mass stopping power	$\text{MeV cm}^2/\text{g}$
Mean excitation potential	eV

7.2 List of particles

Table 7.2: List of available particle identifiers **JPART** and the corresponding particle names.

JPART	Particle	JPART	Particle
-1	All particles	13	Electron
1	Neutron	14	Positron
2	Proton	15	Muon μ^-
3	Pion π^-	16	Muon μ^+
4	Pion π^+	17	e^- -Neutrino ν_e
5	Pion π^0	18	e^- -Anti-neutrino $\bar{\nu}_e$
6	Anti-neutron	19	μ^- -Neutrino ν_μ
7	Anti-proton	20	μ^- -Anti-neutrino $\bar{\nu}_\mu$
8	Kaon κ^-	21	Deuteron
9	Kaon κ^+	22	Triton
10	Kaon κ^0	23	He-3
11	Kaon κ^\sim	24	He-4
12	Gamma ray	25	Heavy ions

Note, the generalized particle **JPART**=-1 is only available within a scoring estimator, and not as a particle source. For a heavy ion **JPART**=25 the charge Z and atomic number A have usually to be defined additionally. As particle source, $Z > 2$ is expected.

7.3 Table of nuclear targets with neutron cross sections

Table 7.3 shows a list of elements and isotopes for which neutron cross sections are included in SHIELD-HIT12A. Natural isotope mixtures are marked with a '*’.

Table 7.3: List of nuclear target identifiers with available neutron cross sections NUCLID and the corresponding names used by SHIELD-HIT12A.

<i>Z</i>	NUCLID	Isotope	Reference
1	1	H-1 - Hydrogen	BNAB-64
1	101	H-2 - Deuterium	BNAB-81
1	102	H-3 - Tritium	BROND
2	104	He-3	BNAB-81
2	2	He-4	BNAB-81
3	105	Li-6	BNAB-81
3	3	Li-7	BNAB-81
4	4	Be-9	BNAB-64
5	106	B-10	BNAB-81
5	5	B-11	BNAB-81
6	6	C-*	BNAB-81
7	7	N-*	BNAB-81
8	8	O-*	BNAB-81
9	9	F-19	BROND-2
11	11	Na-23	BNAB-81
12	12	Mg-*	BNAB-64
13	13	Al-27	BNAB-81
14	14	Si-*	BNAB-81
15	15	P-31	BROND-2
16	16	S-*	JEFF-3
17	17	Cl-*	BROND-2
18	18	Ar-*	JENDL40
19	19	K-*	BNAB-64
20	20	Ca-*	BNAB-81
22	22	Ti-*	JEFF-3
23	23	V-51	BNAB-81
24	24	Cr-*	BNAB-81
25	25	Mn-55	BNAB-81
26	26	Fe-*	BNAB-81
27	27	Co-59	ANL/NDM
28	28	Ni-*	BNAB-81
29	29	Cu-*	BROND-2
30	30	Zn-*	BROND-2
31	31	Ga-*	JENDL
33	33	As-75	JENDL
41	41	Nb-93	BNAB-64
42	42	Mo-*	JENDL

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Table 7.3 – Continued

<i>Z</i>	NUCLID	Isotope	Reference
47	47	Ag-*	JENDL
48	48	Cd-*	BNAB-81
63	63	Eu-*	BNAB-81
64	64	Gd-*	BNAB-81
68	68	Er-*	BNAB-81
73	73	Ta-181	JEFF
74	74	W-*	CENDL31
75	75	Re-*	BNAB-64
79	79	Au-187	BROND-2
80	80	Hg-*	JEFF
82	82	Pb-*	BNAB-81
83	83	Bi-209	BNAB-64
90	90	Th-232	ENDFB-6
92	103	U-235	BNAB-81
92	92	U-238	BNAB-81
94	94	Pu-239	BNAB-81
94	107	Pu-240	BNAB-81

7.4 List of material ICRU_ID

Table 7.4 lists all material numbers as defined by ICRU which are known to SHIELD-HIT12A. The ICRU_ID number can optionally be specified in the material input file *mat.dat* which then will trigger SHIELD-HIT12A to load stopping power data from an external ASCII file (see section 3.2). The file containing the stopping power data *must* exactly be named as the according file name in the table below.

Table 7.4: List of ICRU_ID and file names used by SHIELD-HIT12A.

ICRU_ID	File name	Material name and remarks
1	H.txt	HYDROGEN
2	He.txt	HELIUM
3	Li.txt	LITHIUM
4	Be.txt	BERYLLIUM
5	B.txt	BORON
6	C.txt	AMORPHOUS CARBON (density 2.0 g/cm ³)
906	Graphite.txt	GRAPHITE
7	N.txt	NITROGEN
8	O.txt	OXYGEN
(...)	(...)	(...) ¹
98	Cf.txt	CALIFORNIUM
99	A-150.txt	A-150 TISSUE-EQUIVALENT PLASTIC
100	Acetone.txt	ACETONE
101	Acetylene.txt	ACETYLENE
102	Adenine.txt	ADENINE
103	Adipose.txt	ADIPOSE TISSUE (ICRP)
104	Air.txt	AIR, DRY (NEAR SEA LEVEL)
105	Alanine.txt	ALANINE
106	Al2O3.txt	ALUMINUM OXIDE
107	Amber.txt	AMBER
108	Ammonia.txt	AMMONIA
109	Aniline.txt	ANILINE
110	Anthracene.txt	ANTHRACENE
111	B-100.txt	B100
112	Bakelite.txt	BAKELITE
113	BaF2.txt	BARIUM FLUORIDE
114	BaSO4.txt	BARIUM SULFATE
115	Benzene.txt	BENZENE
116	BeO.txt	BERYLLIUM OXIDE
117	BiGeO.txt	BISMUTH GERMANIUM OXIDE
118	BloodICRP.txt	BLOOD (ICRP)
119	BoneICRU.txt	BONE, COMPACT (ICRU)
120	BoneICRP.txt	BONE, CORTICAL (ICRP)
121	B4C.txt	BORON CARBIDE
122	BoronOxide.txt	BORON OXIDE
123	BrainICRP.txt	BRAIN (ICRP)
124	Butane.txt	BUTANE
125	N-ButylAlcohol.txt	N-BUTYLALCOHOL

Continued on Next Page...

¹All elements up to Californium ($Z = 98$) are supported following this naming scheme.

Table 7.4 – Continued

ICRU_ID	File name	Material name and remarks
126	C-552.txt	C-552 AIR-EQUIVALENT PLASTIC
127	CdTe.txt	CADMIUM TELLURIDE
128	CdWO4.txt	CADMIUM TUNGSTATE
129	CaCO3.txt	CALCIUM CARBONATE
130	CaF2.txt	CALCIUM FLUORIDE
131	CaO.txt	CALCIUM OXIDE
132	CaSO4.txt	CALCIUM SULFATE
133	CaWO4.txt	CALCIUM TUNGSTATE
134	CO2.txt	CARBON DIOXIDE
135	CCl4.txt	CARBON TETRACHLORIDE
136	Cellophane.txt	CELLULOSE ACETATE, CELLOPHANE
137	CAB.txt	CELLULOSE ACETATE BUTYRATE
138	CelNitrate.txt	CELLULOSE NITRATE
139	CeSO4Dos.txt	CERIC SULFATE DOSIMETER SOLUTION
140	CsF.txt	CESIUM FLUORIDE
141	CsI.txt	CESIUM IODIDE
142	ClBenzene.txt	CHLOROBENZENE
143	CHCl3.txt	CHLOROFORM
144	Concrete.txt	CONCRETE PORTLAND
145	Cyclohexane.txt	CYCLOHEXANE
146	DClBenzene.txt	1,2-DICHLOROBENZENE
147	DClDEtyEth.txt	DICHLORODIETHYL ETHER
148	DClEthane.txt	DICHLOROETHANE
149	DEtyEther.txt	DIETHYLETHER
150	DMF.txt	N,N-DIMETHYL FORMAMIDE
151	DMSO.txt	DIMETHYLSULFOXIDE
152	C2H6.txt	ETHANE
153	Ethanol.txt	ETHYL ALCOHOL
154	EthylCell.txt	ETHYL CELLULOSE
155	C2H4.txt	ETHYLENE
156	Eyelens.txt	EYELENS (ICRP)
157	Fe2O3.txt	FERRIC OXIDE
158	FeB.txt	FERRO BORIDE
159	FeO.txt	FERROUS OXIDE
160	FeSO4Dos.txt	FERROUS SULFATE DOSIMETER SOLUTION
161	Freon-12.txt	FREON-12
162	Freon-12B2.txt	FREON-12B2
163	Freon-13.txt	FREON-13
164	Freon-13B1.txt	FREON-13B1
165	Freon-13I1.txt	FREON-13I1
166	GOS.txt	GADOLINIUM OXYSULFIDE
167	GaAs.txt	GALLIUM ARSENIDE
168	GelPhot.txt	GEL IN PHOTOGRAPHIC EMULSION
169	GlassPyrex.txt	GLASS, PYREX
170	GlassLead.txt	GLASS, LEAD
171	GlassPlate.txt	GLASS, PLATE
172	Glucose.txt	GLUCOSE
173	Glutamine.txt	GLUTAMINE
174	Glycerol.txt	GLYCEROL
175	Guanine.txt	GUANINE

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Table 7.4 – Continued

ICRU_ID	File name	Material name and remarks
176	Gypsum.txt	GYPSUM / PLASTER OF PARIS
177	Heptane.txt	N-HEPTANE
178	Hexane.txt	N-HEXANE
179	Kapton.txt	KAPTON POLYIMIDE FILM
180	LaBrO.txt	LANTHANUM OXYBROMIDE
181	La2O2S.txt	LANTHANUM OXYSULFIDE
182	PbO.txt	LEADOXIDE
183	LiNH2.txt	LITHIUM AMIDE
184	Li2CO3.txt	LITHIUM CARBONATE
185	LiF.txt	LITHIUM FLUORIDE
186	LiH.txt	LITHIUM HYDRIDE
187	LiI.txt	LITHIUM IODIDE
188	Li2O.txt	LITHIUM OXIDE
189	Li2B4O7.txt	LITHIUM TETRABORATE
190	Lung.txt	LUNG (ICRP)
191	M3wax.txt	M3 WAX
192	MgCO3.txt	MAGNESIUM CARBONATE
193	MgF2.txt	MAGNESIUM FLUORIDE
194	MgO.txt	MAGNESIUM OXIDE
195	MgB4O7.txt	MAGNESIUM TETRABORATE
196	HgI2.txt	MERCURIC IODIDE
197	CH4.txt	METHANE
198	Methanol.txt	METHANOL
199	MixDwax.txt	MIX D WAX
200	MS20.txt	MS20 TISSUE SUBSTITUTE
201	MuscleICRP.txt	MUSCLE, SKELETAL (ICRP)
202	MuscleICRU.txt	MUSCLE, STRIATED (ICRU)
203	MuscleSuc.txt	MUSCLE EQUIVALENT LIQUID, WITH SUCROSE
204	MuscleNoSuc.txt	MUSCLE EQUIVALENT LIQUID, NO SUCROSE
205	Naphthalene.txt	NAPHTHALENE
206	NitroBenz.txt	NITROBENZENE
207	N2O.txt	NITROUS OXIDE
208	Elvamide.txt	NYLON, DU PONT ELVAMIDE 8062
209	Nylon.txt	NYLON, TYPE 6 AND 6/6
210	Nylon6-10.txt	NYLON, TYPE 6/10
211	Rilsan.txt	NYLON, TYPE 11 (RILSAN)
212	Octane.txt	OCTANE, LIQUID
213	Paraffin.txt	PARAFFINWAX
214	Pentane.txt	N-PENTANE
215	Emulsion.txt	PHOTOGRAPHIC EMULSION
216	Plasscin.txt	PLASTIC SCINTILLATOR (VINYLTOLENE BASED)
217	PuO2.txt	PLUTONIUM DIOXIDE
218	PAN.txt	POLYACRYLONITRILE
219	Polycarb.txt	POLYCARBONATE (MAKROLON, LEXAN)
220	Polyclst.txt	POLYCHLOROSTYRENE
221	Polyeth.txt	POLYETHYLENE
222	Mylar.txt	POLYETHYLENE TEREPHTHALATE (MYLAR)
223	Lucite.txt	POLYMETHYL METHACRALATE (LUCITE, PERSPEX, PMMA)
224	POM.txt	POLYOXYMETHYLENE
225	PP.txt	POLYPROPYLENE

Continued on Next Page...

Table 7.4 – Continued

ICRU_ID	File name	Material name and remarks
226	PS.txt	POLYSTYRENE
227	Teflon.txt	POLYTETRAFLUOROETHYLENE (TEFLON)
228	PTFCE.txt	POLYTRIFLUOROCHLOROETHYLENE
229	PVAc.txt	POLYVINYL ACETATE
230	PVOH.txt	POLYVINYL ALCOHOL
231	PVB.txt	POLYVINYL BUTYRAL
232	PVC.txt	POLYVINYL CHLORIDE
233	Saran.txt	SARAN
234	PVDF.txt	POLYVINYLIDENE FLUORIDE
235	PVP.txt	POLYVINYLPIRROLIDONE
236	KI.txt	POTASSIUM IODIDE
237	K2O.txt	POTASSIUM OXIDE
238	Propane.txt	PROPANE
239	PropaneLiq.txt	PROPANE, LIQUID
240	NPropylOH.txt	N-PROPYL ALCOHOL
241	Pyridine.txt	PYRIDINE
242	RubberB.txt	RUBBER, BUTYL
243	Rubber.txt	RUBBER, NATURAL
244	Neoprene.txt	RUBBER, NEOPRENE
245	SiO2.txt	SILICON DIOXIDE
246	AgBr.txt	SILVER BROMIDE
247	AgCl.txt	SILVER CHLORIDE
248	AgHaEmul.txt	SILVER HALIDES IN PHOTOGRAPHIC EMULSION
249	AgI.txt	SILVER IODIDE
250	Skin.txt	SKIN (ICRP)
251	Na2CO3.txt	SODIUM CARBONATE
252	NaI.txt	SODIUM IODIDE
253	Na2O.txt	SODIUM MONOXIDE
254	NaNO3.txt	SODIUM NITRATE
255	Stilbene.txt	STILBENE
256	Sucrose.txt	SUCROSE
257	Terphenyl.txt	TERPHENYL
258	Testes.txt	TESTES (ICRP)
259	C2Cl4.txt	TETRACHLOROETHYLENE
260	TiCl.txt	THALLIUM CHLORIDE
261	TissueICRP.txt	TISSUE, SOFT (ICRP)
262	TissueICRU.txt	TISSUE, SOFT (ICRU, FOUR COMPONENT)
263	Tissmeth.txt	TISSUE-EQUIVALENT GAS (METHANE BASED)
264	Tissprop.txt	TISSUE-EQUIVALENT GAS (PROPANE BASED)
265	TiO2.txt	TITANIUM DIOXIDE
266	Toluene.txt	TOLUENE
267	Trichlor.txt	TRICHLOROETHYLENE
268	Triethphos.txt	TRIETHYL PHOSPHATE
269	WF6.txt	TUNGSTEN HEXAFLUORIDE
270	UC2.txt	URANIUM DICARBIDE
271	UC.txt	URANIUM MONOCARBIDE
272	UO.txt	URANIUM OXIDE
273	Urea.txt	UREA
274	Valine.txt	VALINE
275	Viton.txt	VITON FLUOROELASTOMER

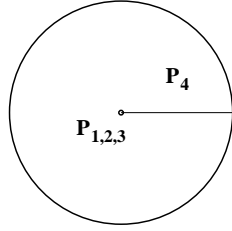
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Table 7.4 – Continued

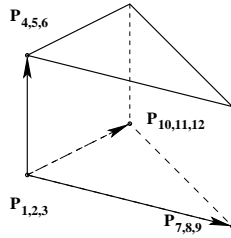
ICRU_ID	File name	Material name and remarks
276	Water.txt	WATER, LIQUID
277	H2Ovapor.txt	WATER VAPOR
278	Xylene.txt	XYLENE

7.5 List of bodies

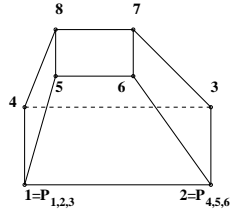
Each body is described with an identifier, that is a three-letter code preceded by two empty spaces. Another integer assigns a number to the body in the next 5 columns. Then 6 blocks each 10 columns wide specify the arguments P_1 to P_6 . If additional arguments are required, a second card is added holding the arguments P_7 to P_{12} , but leaving the first 10 columns empty. Even further cards follow the format of the second card. The arguments P_i can specify a point, a vector or a radius, depending on the body described. The bodies available in SHIELD-HIT12A are listed below:



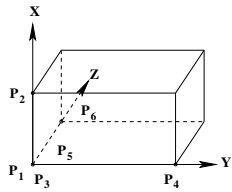
SPH : A sphere defined by the center coordinates (P_1, P_2, P_3) and the radius P_4 . The remaining arguments P_5 and P_6 are ignored.



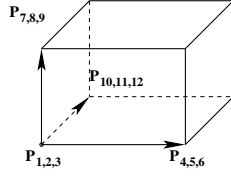
WED : A wedge spanned by three vectors. The origin is the point (P_1, P_2, P_3) from where the three vectors (P_4, P_5, P_6) , (P_7, P_8, P_9) and (P_{10}, P_{11}, P_{12}) start.



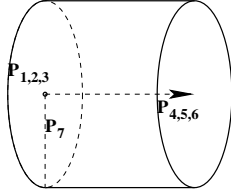
ARB : Arbitrary convex polyhedron. This body is described by 8 points, i.e. 24 arguments. The points must be specified in the order as indicated by the figure to the left: Point 1 is (P_1, P_2, P_3) , point 2 (P_4, P_5, P_6) ... point 8 (P_{22}, P_{23}, P_{24}) .



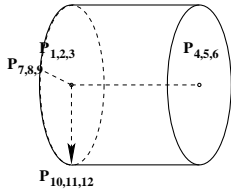
RPP : Rectangular parallelepiped. (P_1, P_2) marks minimum and maximum X coordinates. (P_3, P_4) is min and max values for Y, and (P_5, P_6) for Z.



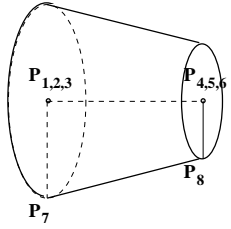
BOX : A box spanned by 3 vectors. All vectors start in the point (P_1, P_2, P_3) and point to the points (P_4, P_5, P_6) , (P_7, P_8, P_9) and (P_{10}, P_{11}, P_{12}) . They must be orthogonal on each other.



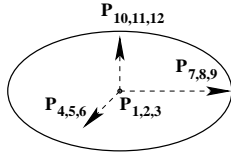
RCC : Right circular cylinder. One end of the cylinder is described by a circle with the center at (P_1, P_2, P_3) . The arguments (P_4, P_5, P_6) hold a vector measured from the center to the opposite end of the cylinder. P_7 marks the radius of the cylinder.



REC : Right elliptical cylinder. Similar to RCC, but instead of a radius, two additional vectors (P_7, P_8, P_9) and (P_{10}, P_{11}, P_{12}) describe the minor and major axis. All three vectors must be orthogonal to each other.



TRC : A cone spanned by two circles. First circle is centered at point (P_1, P_2, P_3) and has radius P_7 . The other end of the cone is described by the circle centered at the end of a vector (P_4, P_5, P_6) starting from the center of the first circle. The second circle has a radius of P_8 . The radius in P_7 must be larger then the radius in P_8 .



ELL : Ellipsoid. Centered at point (P_1, P_2, P_3) , three orthogonal vectors (P_4, P_5, P_6) , (P_7, P_8, P_9) and (P_{10}, P_{11}, P_{12}) span the body.
This body is differently defined than in FLUKA.

MOV : Copies a body number P_1 to a new position translated by the vector (P_2, P_3, P_4) . The columns 6-10 in the MOV card must hold an integer specifying a new and unique body number, just like a normal body specification.

7.6 Input and output files.

Table 7.5 provides an overview of input (I) and output (O) files used by SHIELD-HIT12A at run time. They have either ASCII (A) or binary (B) format. Some files are optional.

Table 7.5: List of input and output files used at run time.

Filename	I/O	ASCII/Bin	Description	Remarks
<i>beam.dat</i>	I	A	Seed, projectile, stats, etc.	
<i>detect.dat</i>	I	A	Auxiliary scoring	1)
<i>detect.bin</i>	O	B	Result of aux. scoring	1), 2)
<i>for017</i>	O	A	GEMCA parser log	
<i>for024</i>	O	A	Broad collection of data	
<i>for025</i>	O	A	Run-time diagnostics	
<i>for026</i>	O	A	Material specific stopping powers, cross sections	
<i>for027</i>	O	A	Neutron cross sections < 14.5 MeV	
<i>for028</i>	O	A	Secondary neutrons production < 14.5 MeV	
<i>for029</i>	O	B	Track length fluence for each zone	
<i>for030</i>	O	B	Double differential particle yields	
<i>for031</i>	O	B	Energy deposition in each zone	
<i>for032</i>	O	B	Energy deposition in zones, diff. in LET	
<i>for033</i>	O	A	Various nuclei distributions in zones	
<i>for034</i>	O	B	Ordinary and planar fluence	
<i>geo.dat</i>	I	A	Geometry description	
<i>mat.dat</i>	I	A	Target medium chemical composition	
<i>parlev.dat</i>	I	A	External parameter file for PARLEV parameters	1), 3)
<i>rifi.dat</i>	I	A	Description of the ripple filter	1), 3)
<i>sobp.dat</i>	I	A	External particle source file	1), 3)

¹⁾ Optional.

²⁾ File name is not predefined but set by user in *detect.dat*.

³⁾ File name is not predefined but set by user in *beam.dat*.

7.7 List of PARLEV parameters

The PARLEV parameters listed in table 7.6 can be used for tuning the nuclear models in SHIELD-HIT. Up to 40 numeric parameters are set in SHIELD-HIT12A which may affect the calculated results. The default values are carefully selected and benchmarked against experimental data, and should therefore not be changed. However, for research purposes they can be changed using an external input file as described in section 3.3.3. In this case, a message is displayed for all modified settings in the output file *for024*.

Values which are marked "for high energies" in the table below, are only used at high energies (well above 1 GeV/A) and will therefore not influence calculations in normal particle radiotherapy settings.

A short description of the individual PARLEV parameters is given below. Details on the individual nuclear fragmentation models are given in the references [13–15, 17–19, 35–38].

PARLEV(1) – PIDABS : The probability of absorption of pion in nucleus by the quasi-deuteron pair.

PARLEV(2) – TLIMIT : Time of simulation of the cascade in nucleus (does not affect the strong decays of the resonances).

PARLEV(3) – MMES : The time of formation of mesons in their rest frame: $\tau = 1/(5.06 \cdot \text{MMES})$ [fm/c].

PARLEV(4) – MBAR : The time of formation of baryons in their rest frame: $\tau = 1/(5.06 \cdot \text{MBAR})$ [fm/c].

PARLEV(5) – EPS1, EPS2 : Separation energy of a nucleon from the nucleus to the cascade stage of the nuclear reaction. The energy is the same for the target nucleus (EPS2) and the projectile (EPS1).

PARLEV(6) – VP : The depth of the potential well for pion inside the nucleus.

PARLEV(7) – C1, C2 : The parameter of diffuseness of the Saxon-Woods distribution of nuclear density $\rho_{\text{WS}}(r) = \rho_0/[1 + \exp(r - R)/c]$, $R = r_0 A^{1/3}$. The parameter is the same for the target nucleus (C2) and the projectile (C1).

PARLEV(8) – D1, D2 : The parameter that determines the maximal radius R_{max} of the distribution of nuclear density $\rho(r)$ from the condition $\rho(R_{\text{max}})/\rho(0) = D$. This parameter is used for both the Saxon-Woods distribution $\rho_{\text{WS}}(r)$ and for light nuclei: $\rho(r) = \rho_0 \exp(-\alpha r^2)$. The parameter is the same for the target nucleus (D2) and the projectile (D1).

PARLEV(9) – RON1, RON2 : The parameter r_0 , which determines the radius of the nucleus by the formula $R = r_0 A^{1/3}$. For light nuclei this parameter directly sets the radius of nucleus. The parameter is the same for the target nucleus (RON2) and the projectile (RON1). PARLEV(9) is applicable to all isotopes, except for those specified in PARLEV(10) – PARLEV(21).

PARLEV(10) – PARLEV(21) : Individual values of RON1, RON2 for light nuclei. SHIELD-HIT12A does not discriminate between RON1 and RON2. In some cases groups of isotopes are affected. For instance, PARLEV(20) affects all nitrogen isotopes with nucleon number of 10 or less. Any isotope which are not affected by PARLEV(10) – PARLEV(21) are treated with PARLEV(9).

- PARLEV(22) - PRECOM : This parameter determines the contribution of pre-equilibrium emission of the lightest nuclei during transition from the cascade stage of nuclear reactions toward the equilibrium de-excitation ($0 \leq \text{PRECOM} \leq 1$).
- PARLEV(23) - ICAN : This parameter determines the number of evaporation channels at equilibrium deexcitation. If ICAN = 0 then only 6 channels are evaporated (n,p,d,t,He³,He⁴+fission). This mode should be used when mainly dealing with Fermi-Breakup of lighter nuclei. If ICAN = 1 then there are 32 channels (+ fission). The latter should only be used when working with heavy excited nuclei.
- PARLEV(24) - IMA : Maximal number of fragments at multifragmentation of nuclei with excitation below 3 MeV/nucleon.
- PARLEV(25) - EPSILO : Inverse level density parameter for internal bulk excitation of fragments in multifragmentation.
- PARLEV(26) - FKACOL : Freeze-out volume for Coulomb interaction of fragments in multifragmentation.
- PARLEV(27) - RNCL : The parameter r_0 , which determines the radius of the excited nucleus by the formula $R = r_0 A^{1/3}$ at the stage of equilibrium de-excitation.
- PARLEV(28) - AMS : Density level parameter of the excited nucleus at evaporation.
- PARLEV(29) - AMFS : Density level parameter of the excited nucleus at fission. Changing this parameter has no effect, as it is overridden internally.
- PARLEV(30) : The border for the Fermi break-up on the number of nucleons: for heavier nuclei the multifragmentation/evaporation/fission model is used.
- PARLEV(31) - ILEVRA : The border for the multifragmentation on the excitation energy per nucleon: for more low excitations only the evaporation/fission model is used. (do not set higher than 2 MeV/nucleon).
- PARLEV(32) - ILEVRA : Products of Fermi break-up are produced only in the ground state (ILEVRA=1), or these products can be formed in an excited state (ILEVRA=2).
- PARLEV(33) - FKAP1 : Fermi break-up: freeze-out volume for translation motion of fragments.
- PARLEV(34) - FKAP2 : Fermi break-up: freeze-out volume for Coulomb interaction of fragments.
- PARLEV(35) -- PARLEV(38) : Reserved.
- PARLEV(39) - SIGION : The parameter for renormalization the total $\sigma_{\text{tot}}(E)$ and inelastic $\sigma_{\text{in}}(E)$ cross sections of nucleus-nucleus interactions in the cascade inside a macroscopic target.
- PARLEV(40) - MICROD : The parameter for renormalization the total $\sigma_{\text{tot}}(E)$ and inelastic $\sigma_{\text{in}}(E)$ cross sections of hadron-nucleus interactions in the cascade inside a macroscopic target.

Table 7.6: List of PARLEV parameters with their default values in SHIELD-HIT12A.

PARLEV#	Description	Default	Units	Remarks
1	AGT: PIDABS	0.01	probability	
2	AGT: TLIMIT	25.0	fm / c	¹⁾
3	AGT: MMES	0.2	GeV	¹⁾
4	AGT: MBAR	999.999	GeV	¹⁾
5	AGT: EPS1,EPS2	0.007	GeV	
6	AGT: VPI	0.025	GeV	
7	AGT: C1,C2	0.545	fm	Wood-Saxon param.
8	AGT: D1,D2	0.05	(N/A)	
9	AGT: R0N1,R0N2	1.097	fm	
10	AGT: R0N H(2,1)	2.704	fm	
11	AGT: R0N H(3,1)	2.55	fm	
12	AGT: R0N He(A,2)	2.55	fm	$A \neq 4$
13	AGT: R0N He(4,2)	2.157	fm	
14	AGT: R0N Li(6,3)	3.317	fm	
15	AGT: R0N Li(A,3)	3.11	fm	$A \neq 6$
16	AGT: R0N Be(A,4)	3.25	fm	all Be isotopes
17	AGT: R0N B(10,5)	3.16	fm	
18	AGT: R0N B(A,5)	2.48	fm	$A < 10$
19	AGT: R0N C(A,6)	2.48	fm	$A \leq 10$
20	AGT: R0N N(A,7)	2.48	fm	$A \leq 10$
21	AGT: R0N O(A,8)	2.48	fm	$A \leq 10$
22	PRECO: PRECOM	1.0	(N/A)	
23	DEEX: ICAN	0.0	flag: 0 or 1	
24	DEEX: IMA	3.0	(N/A)	
25	DEEX: EPSIL0	16.0	MeV	
26	DEEX: FKACOL	2.0	(N/A)	
27	DEEX: RNCL	0.0	fm	
28	DEEX: AMS	0.125	MeV ⁻¹	level density
29	DEEX: AMFS	0.125	MeV ⁻¹	level density
30	DEEX: PARLEV(30)	16.0	Mass number	A_{\max} for Fermi density
31	DEEX: PARLEV(31)	0.002	GeV/nucleon	min of U_{mult}/A
32	DEEX: ILEVRA	2.0	flag: 1 or 2	ground or excited
33	DEEX: FKAP1-Coulomb	0.65	(N/A)	
34	DEEX: FKAP2-Volume	18.0	(N/A)	
35	(Reserved)	0.0	-	
36	(Reserved)	0.0	-	
37	(Reserved)	0.0	-	
38	(Reserved)	0.0	-	
39	SIGION	1.00	(N/A)	Renorm $\sigma_{AA}(E)$
40	MICROD	1.00	(N/A)	Renorm $\sigma_{hA}(E)$

¹⁾ For high energies1 fm = 10^{-13} cmc = $3 \cdot 10^{10}$ cm/sec

N/A : Non-applicable, dimensionless

Chapter 8

Legacy and additional background information

8.1 The legacy format of target medium file *mat.dat*

Usage of the old format for *mat.dat* is discouraged, instead users should migrate to the new format as described in section 3.2.

In the old format, four different types MEDTYP of media are considered in the SHIELD-HIT code:

1. Pure medium MEDTYP=1 - element with fixed Z (e.g. O, Fe, Cu etc.)
2. Chemical composition MEDTYP=2, e.g. water H_2O .
3. Isotope mixture MEDTYP=3, if isotopes of the same Z have different neutron properties at low energies, e.g. mixture of ^{235}U and ^{238}U .
4. Others MEDTYP=4, which do not conform to types 1-3. In this case the user should calculate the atomic concentrations (in units 10^{27} atoms/cm³) and the partial densities of elements (in units g/cm³) manually.

The total number of different media NUMMED in any task may be up to 16, excluding internal and outer vacuum which have the values 1000 and 0, respectively.

In order to define a MEDTYP the user has to specify all chemical elements which can be found in this medium. A chemical element is described by the variable NUCLID. Usually $\text{NUCLID}=Z$ selects the element from the periodic table corresponding to the atomic number Z ($1 < Z < 100$). Additionally, there are several isotopes with a different number of neutrons for which a specific value of $\text{NUCLID} > 100$ is already predefined. A list of these predefined values is given in table 7.3 in section 7.2. The number NELEM of elements (i.e. of variables NUCLID) in each medium may be up to 13.

Several examples of *mat.dat* are given below. *It has to be emphasized that all symbol positions in the files mat.dat (excluding comment lines) must be strictly kept!*

MEDTYP=1: pure medium

This is the most basic MEDTYP. Only one NUCLID per medium is expected.

To define a chemical composition, the user should set up the variable **NUCLID** for each element and numeric indexes of the chemical formula. Moreover, the user should define the number of constituent elements in the variable **NELEM** and the total density of the composition **RHO** (in g/cm^3).

Optionally, the mean excitation energy I (eV) may be added at the very right (column 57 and beyond) of the line to overwrite the default values. Otherwise, the optional variable **P** which follows as fourth argument after **RHO** can take the values 0, 1, and 2. It can be used to invoke the I values as recommended by ICRU [27], that is,

P=1 : I values for elements forming a gas

P=2 : I values for elements in a liquid or solid compound

[illegible]

To define a mixture of isotopes (MEDTYP=3) the user should, as for MEDTYP=2, define the variables **NELEM** and **RHO** specifying the number of elements and the total density (in g/cm³) of the mixture. Moreover, the user should set up the variables **NUCLID** for each isotope and **PART** for the relative fraction of each isotope of the number of atoms. For example, natural uranium is defined this way:


```

NUMMED= 1
                                MEDIUM NO. 1 (NatU)
MEDTYP=3 NELEM=2 RHO=18.900
1 NUCLID=103 PART=0.70000E-02
2 NUCLID= 92 PART=0.99300E 00

```

MEDTYP=4: arbitrary mixture of elements

The same medium which is defined in the example for MEDTYP=3, i.e. natural uranium, can also be obtained by using the medium of type 4:

```

NUMMED= 1
                                MEDIUM NO. 1 (NatU)
MEDTYP=4 NELEM 2 RHO=18.900
1 NUCLID=103 DEL1=.334740E-06 RH01= .130628
2 NUCLID= 92 DEL2=.474853E-04 RH02=18.769372

```

To define an arbitrary mixture, the user should set up the variable NUCLID for each element, partial concentrations, and partial densities for each NUCLID. Like above, the user should define also the variable NELEM and the total density of the mixture RHO (in g/cm³). Moreover, the partial concentrations for each NUCLID in units 10²⁷ atoms/cm³ and the partial densities for each NUCLID in units g/cm³, should be given.

```

NUMMED= 1
                                MEDIUM NO. 1 (Air: N 0.75527, O 0.2318, w/o Ar)
ICRU_ID=104
MEDTYP=4 NELEM 3 RHO=0.001203 P=1
1 NUCLID= 6 DEL1=1.48863E-11 RH01=1.49395E-07
2 NUCLID= 7 DEL2=3.93565E-08 RH02=9.22040E-04
3 NUCLID= 8 DEL3=1.05672E-08 RH03=2.82800E-04

```

The ICRU_ID can be used to choose stopping power data from an external file as described in Sec. 3.1 above and Table 7.4 lists the ICRU_ID and naming conventions. If ICRU_ID is set to 0 or the specified file is missing, then SHIELD-HIT12A will use its own default stopping power data which are precalculated using an internal modified Bethe-Bloch with Linhard-Scharff formula and will additionally output a warning in the latter case.

External stopping power files containing mass stopping power data for a set of materials and for ions from ¹H⁺ (proton) to ⁴⁰Ar¹⁸⁺ can be provided by the user if the ICRU_ID flag is set to an appropriate value from the list in table 7.4. These external mass stopping power files are expected to be ASCII formatted, space or tab separated columns, and should follow the ICRU 73 [25] energy grid (53 energy nodes from 25 keV/u up to 1 GeV/u). SHIELD-HIT12A expects the grid to contain the first 18 ions, i.e. from protons ($Z=1$) to Argon ions ($Z=18$). The mass stopping power data must be in units of MeV cm²/g. Comment lines starting with * are allowed anywhere in the file, but no line (either comment or data) may be longer than 512 characters. External stopping power files may contain data different from ICRU, e.g. ATIMA [2] data, but the energy grid must be compatible to ICRU 73. The libdEdx [3, 26] library can provide a source for stopping power data. A script *shield_dEdx* described in section 5.2.4 can generate these tables for most ICRU materials using libdEdx.

Examples of external stopping power files are: *Water.txt*, *Air.txt*, *A-150.txt*, *Kapton.txt*, which are supplied along with the SHIELD-HIT12A distribution. In table 7.4 the naming convention of the external stopping power files is listed.

8.2 History

This is a short ChangeLog that describes the most noteworthy changes in the A-branch of SHIELD-HIT.

SHIELD-HIT12A

SHIELD-HIT12A is the first version of SHIELD-HIT that was ever released in public.

- All changes from SHIELD-HIT10 were merged into SHIELD-HIT12A; these were mostly minor bugfixes.
- *for022.dat*, *for023.dat* and *pasin.dat* were renamed into *mat.dat*, *beam.dat* and *geo.dat*
- Comments allowed in *geo.dat*.
- *mat.dat* is now in free format, making extension with new cards possible without breaking compatibility. Comments are now allowed, too.
- *beam.dat* is now in free format, making extension with new cards possible without breaking compatibility. Comments are now allowed, too.
- Flat circular and flat square beams are now possible.
- Transportation in any direction, specified with the **BEAMDIR** card.
- Implemented Gaussian or flat optional beam divergence and beam focus model using the **BEAMDIV** card.
- Updated inelastic antiproton cross sections. This is still experimental.
- Implementation of a new Vavilov straggling function, which is 5-6 times faster than the old version taken from GEANT3.21. This speeds up SHIELD-HIT12A by 30-40 % for a typical C-12 depth dose calculation.
- New Moliere scattering function.
- Removed code that was taken from GEANT3.21 thereby liberating SHIELD-HIT12A from GPL license and Copyright issues.
- No need for *atab.dat* and *tabnuc.dat* files as they are hardcoded now.
- Bug fixes, most prominently the missing fluence in vacuum issue (ticket #27).
- Script for generating stopping power data files from libdEdx.
- Scoring by zone is now provided by the new scoring system. Zone-scoring was also possible with the old scorer, however, incompatible with parallelization.
- New detector: alanine dosimeter response model is now included following [32].
- Implementation of neutron cross sections (below 14.5 MeV) for Argon. Update of neutron cross sections for Tungsten.

SHIELD-HIT10A

SHIELD-HIT10A was never released as such, but was subject to continuous development. SHIELD-HIT10A started as a fork from the SHIELD-HIT08 code. Changes to SHIELD-HIT10 were still migrated to SHIELD-HIT10A though during the development process.

- This manual was initiated.
- The *shieldhit* executable accepts a directory as argument, and -n option provides an iterator for multiple runs in the same directory, which is useful for parallelization.
- Arbitrary mesh / cylindrical scoring, with several estimators.
- External beam files can be read, i.e. Spread out Bragg-peak data files derived from GSI raster scan files.
- Ripple filters can be defined and loaded.
- RANLUX random number generator replaced by RANSHI.
- Added script *shield_detect2ascii* which can average results from multiple parallelized runs.
- Baraschenkov's cross sections updated for heavy ion - heavy ion reactions.
- Fermi-Breakup: free coulomb energy and free volume parameters changed from 1.0 to 0.65 and 18.0 respectively (see PARLEV33 and 34).

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