

Computation Tools for High Energy Physics and cosmology.

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MicrOMEGAs, A code for calculation of Dark Matter signals in Beyond Standard Models

<https://lapth.cnrs.fr/micromegas>

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MicrOMEGAs team

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General features and purpose

- MicrOMEGAs is based on CalcHEP package

<https://theory.sinp.msu.ru/~pukhov/calchep.html>

which is intended for calculation of cross sections and particle decay widths in generic model of particle interaction.

- MicrOMEGAs is able to calculate relic density of DM, signals of direct and indirect DM detection.
- Micromegas contains/imports external packages for model construction, calculation of particle spectra, testing of collider signals.
- Operation system Linux or Darwin (Mac)
- User's code language is C or C++

Downloading of micrOMEGAs package

micrOMEGAs site

<http://lapth.in2p3.fr/micromegas>

Click *Download and Install* (left -top part of the screen)

And then *DOWNLOAD* (right-top part of the screen)

The name of received file should be

micromegas_5.2.10.tgz

Unpack it by **tar -xvzf micromegas_5.2.10.tgz**

It should create directory **micromegas_5.2.10/** which occupies about 180 Mb of disk space. You will need more disk space after compilation of specific models and generation of matrix elements .

Compilation of micromegas code consists of 3 steps.

- 1) Compilation of CalcHEP and main micrOMEGAs routines
- 2) Compilation of code for given model of particle interaction.
- _3) Runtime compilation of external packages and matrix elements.

File structure of micrOMEGAs package.

micromegas_5.2.10/	main directory
CalcHEP_src/	generator of matrix elements
sources/	micrOMEGAs own codes
man/	description of micrOMEGAs routines manual_5.2.pdf
include/	micromegas.h & micromegas_aux.h
lib/	for generated libraries
Packages/	
	SuSpect_2.41 NMSSMTools_4.7.1 CpsuperH2.3, LoopTools-2.1 LanHEP
<i>model directories:</i>	
MSSM/	
NMSSM/	Next-to-Minimal SuSy Model
CPVMSSM/	MSSM with complex parameters
IDM/	Inert doublet model
LHM/	Little Higgs Model
Z3IDM/	Inert doublet and singlet, Z^3 symmetry
Z4IDSM/	Inert doublet and singlet, Z^4 symmetry

Structure of MODEL directory

main.c[pp]	file with <i>main</i> program for given model
Makefile	supports compilation of C and C++ user codes
lib/	
*.c, .F,cpp	source codes of specific model routines
Makefile	called automatically to generate
alib.a	compiled library
work/	CalcHEP working directory intended for matrix element generation
models/	model in CALCHEP format:
	vars1.mdl func1.mdl prtcls1.mdl
	lgrng1.mdl
so_generated/	directory to store automatically generated matrix elements
./calchep	launch interactive CalcHEP session

Structure of main.c file

main.c, main.cpp files in micrOMEGAs model directories consist of several independent blocks enclosed into

```
#ifdef XXXXX
```

```
.....
```

```
#endif
```

In the top of main.c the user can switch on/off any of this block via corresponding *#define* instruction at the top of file

```
#define MASSES_INFO           // Display information about mass spectrum
#define CONSTRAINTS          // Display B->s,gamma, Bs->mu,mu,
#define LILITH                // Test of Higgs properties
#define HIGGSBOUNDS
#define OMEGA                 // Calculate relic density
#define INDIRECT_DETECTION   // Signals of DM annihilation in galaxy halo
//#define RESET_FORMFACTORS  // Redefinition of Form Factors and other parameters
#define CDM_NUCLEON          // Calculate amplitudes and cross-sections for CDM-
                             // nucleon collisions
//#define CDM_NUCLEUS        // Calculate number of events for 1kg*day and recoil
                             // energy distribution for various nuclei
#define NEUTRINO              // neutrino telescope
#define DECAY                 // particle width and decay branching
#define CROSS_SECTIONS       // calculate cross sections
```

The main.c files from all model directories are similar and call the same micrOMEGAs routines.

Special requirement for compilation

- a) **gfortran**. Mac users has to pay attention on C-Fortran compatibility.
- b) **X11 develop package** (file /usr/include/X11/*.h):
Needed for Graphic User Interface in CalcHEP and plot facilities of micrOMEGAs. One can **ignore** this requirement or install
- libX11-devel for Fedora/Scientific, old Mac
 - Xquartz (<https://www.xquartz.org>) , new Mac
 - libX11-dev for Ubuntu/Debian [old Ubuntu]
 - libx11-dev for Ubuntu/Debian [new Ubuntu]
 - xorg-x11-devel for SUSE
- c) SMODELS external package requires **python 3.0** or later

Compilation of micrOMEGAs routines

To compile CalcHEP and main micrOMEGAs routines one has to call

make

In micrOMEGAS directory. Compiler flags are stored here

`CalcHEP_src/FlagsForSh`

Any model directory contains `main.c` & `Makefile`

Command

make

generates executable

`./main`

In general

make main=*my_main.c*

generates executable file

`./my_main`

Default `main.c` codes generate executable which needs one argument, the name of file with input parameters. For instance

`./main data1.par`

Run time compilation of matrix elements

If micrOMEGAs needs matrix element for some process, or structure of model vertex, it calls CalcHEP for matrix element generation. Code of matrix element/vertex is compiled, presented as a shared library and stored in directory `MODEL/work/so_generated`.

User sees the message on the screen

```
PROCESS: <name of process>
```

Or

```
VERTEX : < name of vertex>
```

Information about vertices is used to compile effective loop induced Higgs-photon and Higgs-gluon vertices.

Shared library is loaded dynamically in run time

Each shared libraries generated only one time.

If model of interaction is changed, then shared library is recompiled automatically

Dark Matter in micrOMEGAs models.

Discrete symmetry.

MicrOMEGAs assumes a discrete symmetry which is responsible for stability of Dark Matter. For instance, it could be a Z_2 symmetry which divides all particles in two classes, **odd** and **even**, say R-parity in MSSM. The **lightest odd** particle is stable and can be treated as DM.

For micrOMEGAs odd particles are particles whose name started with tilde “~”. For example, ~X, ~H3, ~H+ in IDM.

In case of Z_4 symmetry internal charge for DM particles can be +/- 1 or 2. DM1- the lightest particle with charge 1 is always stable. But the lightest particle with charge 2 is stable if its mass is less than the mass of 2 DM1 particles. One can also construct a model with complex symmetry like $Z_2 \times Z_3$ which always has 2 DM particles.

MicrOMEGAs can work with models with 2DM classes which are marked by “~” and “~~”

Example: Inert Doublet Model

Inert Doublet model contains two $SU(2) \times U(1)$ doublets

$$H_1 = \begin{pmatrix} 0 \\ \langle v \rangle + h/\sqrt{2} \end{pmatrix}, \quad H_2 = \begin{pmatrix} \tilde{H}^+ \\ (\tilde{X} + i \cdot \tilde{H}^3)/\sqrt{2} \end{pmatrix}$$

The Lagrangian contains only even powers of H_2 doublet

$$L = (SM \text{ terms}) + D^\mu H_2^* D_\mu H_2$$

$$-\mu^2 H_2^2 - \lambda_2 H_2^4 - \lambda_3 H_1^2 H_2^2 - \lambda_4 |H_1^* H_2|^2 - \lambda_5 \text{Re}[(H_1^* H_2)^2]$$

Because of symmetry $H_2 \rightarrow -H_2$ the lightest of $\tilde{H}^+, \tilde{X}, \tilde{H}^3$ is stable

Parameters $\mu, \lambda_3, \lambda_4$ can be expressed in terms of masses

New couplings are $\lambda_2, \lambda_L = \lambda_3 + \lambda_4 + \lambda_5$

See details in [arXiv:1106.1719](https://arxiv.org/abs/1106.1719)

vars1.mdl: Free parameters of the model.

Inert Doublet Model
Variables

Name	Value	> Comment	<
EE	0.31333	Electromagnetic coupling constant	
SW	0.474	sin of the Weinberg angle	
MZ	91.187	Mass of Z	
MHX	111	Mass of Inert Doublet Higgs	
MH3	222	Mass of CP-odd Higgs	
MHC	333	Mass of charged Higgs	
LaL	0.01	Coupling in Inert Sector	

.....

func1.mdl: Constrained parameter of the model.

Inert Doublet

Constraints

Name	> Expression
CW	sqrt(1-SW^2)
MW	MZ*CW
Mb	MbEff(Q)
Mc	McEff(Q)
mu2	MHX^2 - laL*(2*MW/EE*SW)^2
la3	2*(MHC^2 - mu2)/(2*MW/EE*SW)^2
la5	(MHX^2 - MH3^2)/(2*MW/EE*SW)^2

prtcls1.mdl: Particles of the model

List fo particles presented in file MODEL/work/models/prtcls1.mdl

Full Name	P	aP	number	spin2	mass	width	color	aux	> LaTeX(A)
photon	A	A	22	2	0	0	1	G	A
Z boson	Z	Z	23	2	MZ	!wZ	1	G	Z
gluon	G	G	21	2	0	0	8	G	G
W boson	W+	W-	24	2	MW	!wW	1	G	W ⁺
neutrino	n1	N1	12	1	0	0	1	L	\nu ^e
electron	e1	E1	11	1	0	0	1		e
mu-neutrino	n2	N2	14	1	0	0	1	L	\nu ^{\mu}
muon	e2	E2	13	1	Mm	0	1		\mu
tau-neutrino	n3	N3	16	1	0	0	1	L	\nu ^{\tau}
tau-lepton	e3	E3	15	1	Mt	0	1		\tau
u-quark	u	U	2	1	0	0	3		u
d-quark	d	D	1	1	0	0	3		d
c-quark	c	C	4	1	Mc	0	3		c
s-quark	s	S	3	1	Ms	0	3		s
t-quark	t	T	6	1	Mtop	wtop	3		t
b-quark	b	B	5	1	Mb	0	3		b
Higgs	h	h	25	0	Mh	!wh	1		h
odd Higgs	~H3	~H3	36	0	MH3	!wH3	1		(H3)
Charged Higgs	~H+	~H-	37	0	MHC	!wHC	1		(H+)
second Higgs	~X	~X	35	0	MHX	!wHX	1		(X)

Names of particles of **odd** sector start with tilde ~

lg RNG1.mdl: Feynman rules

Inert Doublet
Lagrangian

P1	P2	P3	P4	> Factor	< > dLagrangian/ dA(p1) dA(p2)dA(p3)
A	W+	W-		-EE	m3.p2*m1.m2 - m1.p2*m2.m3 -
A	~H+	~H-		EE	m1.p3 - m1.p2
B	b	A		EE/3	G(m3)
B	b	G		GG	G(m3)
B	b	Z		-EE/(12*CW*SW)	4*SW^2*G(m3) - 3*G(m3)*(1-G5)
B	b	h		-EE*Mb/(2*MW*SW)	1
B	t	W-		-EE*Sqrt2/(4*SW)	G(m3)*(1-G5)
W+	W-	~X	~X	EE^2/(2*SW^2)	m1.m2
h	~X	~X		-2*MW*SW/EE	la3+la4+la5
Z	Z	~X	~X	EE^2/(2*CW2*SW^2)	m1.m2

.....

p – momentum, **m** – Lorentz index

Example of micrOMEGAs session for IDM

./main data1.par

VERTEX: W- W+ h
VERTEX: L l h
VERTEX: C c h
VERTEX: T t h
VERTEX: B b h
VERTEX: ~H- ~H+ h

Dark matter candidate is '~X' with spin=0/2

=== MASSES OF HIGGS AND ODD PARTICLES: ===

Higgs masses and widths

PROCESS: h->2*x

PROCESS: W+->2*x

PROCESS: Z->2*x

PROCESS: h->W-,E,ne

Delete diagrams with W+<1

PROCESS: h->Z,ne,Ne

Delete diagrams with Z<1

h 125.00 3.97E-03

Masses of odd sector Particles:

~X : MHX = 600.0 || ~H3 : MH3 = 601.0 || ~H+ : MHC = 604.0

LILITH(DB15.09): -2*log(L): 25.96; -2*log(L_reference): 0.00; ndf: 38; p-value: 9.31E-01

Continue

==== Calculation of relic density =====

PROCESS: $\sim X, \sim X \rightarrow \text{AllEven}, 1 * x \{A, Z, G, W+, W-, ne, Ne, e, E, nm, Nm, m, M, nl, NI, l, L, u, U, \dots\}$

PROCESS: $\sim H3, \sim X \rightarrow \text{AllEven}, 1 * x \{A, Z, G, W+, W-, ne, Ne, e, E, nm, Nm, m, M, nl, NI, l, L, u, U, \dots\}$

PROCESS: $\sim H3, \sim H3 \rightarrow \text{AllEven}, 1 * x \{A, Z, G, W+, W-, ne, Ne, e, E, nm, Nm, m, M, nl, NI, l, L, \dots\}$

.....

Xf=2.62e+01 Omega=1.13e-01

Channels which contribute to 1/(omega) more than 1%.

Relative contributions in % are displayed

21% $\sim X \sim X \rightarrow W+ W-$

14% $\sim X \sim X \rightarrow Z Z$

11% $\sim H3 \sim H3 \rightarrow W+ W-$

9% $\sim H+ \sim H- \rightarrow W+ W-$

7% $\sim H3 \sim H3 \rightarrow Z Z$

6% $\sim H+ \sim X \rightarrow A W+$

5% $\sim H3 \sim H+ \rightarrow A W+$

4% $\sim H+ \sim H- \rightarrow A A$

4% $\sim H3 \sim H+ \rightarrow Z W+$

3% $\sim H+ \sim X \rightarrow Z W+$

3% $\sim H+ \sim H- \rightarrow A Z$

2% $\sim H+ \sim H- \rightarrow Z Z$

2% $\sim H+ \sim X \rightarrow W+ h$

1% $\sim H+ \sim H- \rightarrow h h$

==== Indirect detection =====

annihilation cross section $6.18E-26 \text{ cm}^3/\text{s}$

contribution of processes

$\sim X, \sim X \rightarrow W^+ W^-$ $6.01E-01$

$\sim X, \sim X \rightarrow Z Z$ $3.99E-01$

$\sigma_{\text{av}} = 6.18E-26 [\text{cm}^3/\text{s}]$

Photon flux for angle of sight $f = 0.10 [\text{rad}]$

and spherical region described by cone with angle $0.10 [\text{rad}]$

Photon flux = $9.37E-16 [\text{cm}^2 \text{ s GeV}]^{-1}$ for $E = 300.0 [\text{GeV}]$

Positron flux = $1.04E-13 [\text{cm}^2 \text{ sr s GeV}]^{-1}$ for $E = 300.0 [\text{GeV}]$

Antiproton flux = $5.91E-13 [\text{cm}^2 \text{ sr s GeV}]^{-1}$ for $E = 300.0 [\text{GeV}]$

==== Calculation of CDM-nucleons amplitudes =====

PROCESS: QUARKS, $\sim X \rightarrow$ QUARKS, $\sim X\{u, U, d, D, c, C, s, S, t, T, b, B$

Delete diagrams with $_S0_ \neq 1, _V5_ , A$

CDM[antiCDM]-nucleon micrOMEGAs amplitudes:

proton: SI $1.497E-11 [1.497E-11]$ SD $0.000E+00 [0.000E+00]$

neutron: SI $1.512E-11 [1.512E-11]$ SD $0.000E+00 [0.000E+00]$

CDM[antiCDM]-nucleon cross sections[pb]:

proton SI $9.767E-14 [9.767E-14]$ SD $0.000E+00 [0.000E+00]$

neutron SI $9.962E-14 [9.962E-14]$ SD $0.000E+00 [0.000E+00]$

=====Neutrino Telescope===== for Sun

$E > 1.0E+00 \text{ GeV}$ neutrino/anti-neutrino fluxes $1.81E+01/2.05E+01 [1/\text{Year}/\text{km}^2]$

IceCube22 exclusion confidence level = $1.29E-07\%$

$E > 1.0E+00 \text{ GeV}$ Upward muon flux $2.337E-07 [1/\text{Year}/\text{km}^2]$

$E > 1.0E+00 \text{ GeV}$ Contained muon flux $6.999E-07 [1/\text{Year}/\text{km}^3]$

Main functions of micrOMEGAs

Free model parameters.

assignValW(*name,value*) assigns new value to parameter.

For example

```
assignValW("MHX",600)
```

Function to download set of parameters:

```
readVar(fileName)
```

Structure of file has to be

```
name      value    [ # comment ]
```

For instance, in case of IDM

```
1aL      0.001    # coupling
MHX      600     # inert sector Higgs
Mh       125    # SM Higgs mass
1a2      0.01   # coupling
MHC      604    # mass of charged Higgs
MH3      601    # mass of CP odd Higgs
```

Checking of constrained models parameters

After assignment of parameters one has to call

sortOddParticles(outText);

which calculates constrained parameters and finds DM particle[s]. In case of error in calculation of constrained parameter this routine returns error code and *outText* contains the name of parameter which can not be calculated.

In case of success **sortOddParticles** detects the lightest odd particle[s]

CDM1 [CDM2] and their masses

Mcdm1 [Mcdm2, Mcdm=min(Mcdm1,Mcdm2)]

Values of constrained parameters can be obtained by

findValW(*name*)

Masses of particles can be obtained by

pMass(*name*)

Calculation of DM relic density for 1 DM case

darkOmega(&Xf,fast,Beps)

uses Runge-Kutta to solve

$$\frac{dY}{ds} = \frac{\langle v\sigma \rangle}{3H} (Y^2 - Y_{eq}^2)$$

Returns Ωh^2

fast = 1 for fast calculation: Gauss n-point integration with separation of s-channel poles

0 for precise calculation, Simpson adaptive integration

-1 for very precise Simpson adaptive integration with separation of s-channel poles

Beps removes co-annihilation if

$$\exp\left(\frac{2M_{cdm} - M_1 - M_2}{T}\right) < Beps$$

$X_f = M_{cdm}/T_f$, where $Y(T_f) = 2.5 Y_{eq}(T_f)$ defines freeze-out temperature

printChannels(Xf, Beps, cut,prc,file) print out main annihilation channels and their contributions to $1/\Omega h^2$

21% $\sim X \sim X \rightarrow W^+ W^-$

14% $\sim X \sim X \rightarrow Z Z$

11% $\sim H^3 \sim H^3 \rightarrow W^+ W^-$

9% $\sim H^+ \sim H^- \rightarrow W^+ W^-$

Direct Detection

To predict results of direct detection experiment in a given model we have to calculate cross sections of DM – nuclei elastic scattering. Velocities of DM particles in halo of Milky Way are about orbital velocities of stars $\sim 0.001c$. So, we calculate elastic cross sections in $v=0$ limit.

Feynman rules give us

DM - quarks cross section

Then using nucleon form factors we get

DM - nucleon scattering cross section

Then using nucleus form factors we get

DM -nucleus cross section

In the last step we perform recasting of Direct Detection experiments to make conclusion about validity of the model.

There are Spin Independent (scalar) and Spin Dependent (spin flip) cross section. Even (particle \leftrightarrow antiparticle) and odd cross sections.

Nucleon amplitudes and cross sections in micrOMEGAs

nucleonAmplitudes(name_of_DM ,pA0,pA5,nA0,nA5);

Output: *pA0,pA5,nA0,nA5* – 2 dimension arrays

Proton

pA0[even SI, odd SI] pA5[even SD, odd SD]

Neutron

nA0[even SI, odd SI] nA5[even SD, odd SD]

Then DM-nucleon cross section in [pb] units are

$$\sigma_{SI} = C * A^2 \quad \sigma_{SD} = 3 * C * A^2 \quad \text{where } C = 4/\pi * 3.89E8 * (M_N * M_{dm} / (M_N + M_{dm}))^2$$

CDM[antiCDM]-nucleon micrOMEGAs amplitudes:

proton: SI 1.497E-11 [1.497E-11] SD 0.000E+00 [0.000E+00]

neutron: SI 1.512E-11 [1.512E-11] SD 0.000E+00 [0.000E+00]

micrOMEGAs function for nuclei

nucleusRecoil(

f, - velocity distribution $f(v[\text{km/s}])$ normalized by

$$\int_0^{\infty} f(v)dv = 1$$

A, - atomic number

Z, - nucleus charge

J, - number of spin states

Sxx, - SD formfactors

dNdE - recoil energy distribution stored in array

)

dNdERecoil(E[keV],dNdE) interpolates dNdE table and gives spectrum in 1/keV/kg/day units

For example:

```
nEvents=nucleusRecoil(Maxwell,73,Z_Ge,J_Ge73,SxxGe73,dNdE);
```

Result depends on **global parameter**

rhoDM 0.3[GeV/cm³] Dark Matter density at R_{sun}

Vrot 220[km/s] Galaxy rotation velocity at R_{sun}

Vearth 225[km/s] Galaxy velocity of the Earth

Vesc Escape velocity at R_{sun}

Comparison with DD experiments

$\alpha = 1\text{-C.L.}$ can be obtained by

`DD_pval(expCode, $f\nu$, &expName)`

expCode can include

`XENON1T_2018, DarkSide_2018, PICO_2019, CRSST_2019`

For example

`allExp=XENON1T_2018 | DarkSide_2018 | PICO_2019 | CRSST_2019`

expName – the name of experiment that leads to the strongest exclusion

arXiv:0803.2360 [hep-ph]

arXiv:2003.08621 [hep-ph]

Indirect detection in micrOMEGAs

Indirect detection -detection of photons, positrons and antiprotons signal obtained in result of DM annihilation in Galactic Halo.

For various spectra we use $NZ=250$ dimension arrays and interpolation function for them is `SpectdNdE(E,spectArr)`

One can use `displayPlot` to see and compare different spectra.

`vsigma=calcSpectrum(key,Sg,Se,Sp,Sne,Snm,Snl,&err)`

Calculates $v\sigma$ cross section in cm^3/sec units for DM annihilation photon `Sg`, positron `Se`, antiprotons `Sp`, and neutrino `Sne,Snm,Snl` contain spectra for DM-DM annihilation

Here the average over DM,DM/antiDM is done. PYTHIA 6.4 was used for hadronisation of primary annihilation channels.

Meaning of **key parameter**:
1-takes into account W/Z polarization
2-include gammas from $2 \rightarrow 2 + \text{gamma}$
4-print cross sections

loopGamma

For all implemented models we have

$DM, DM \rightarrow \text{photon}, \text{photon}$ and $DM, DM \rightarrow \text{photon}, Z$ loop induced signals. These signals are not compiled automatically in run-time but generated in advance by means of `FormCalc`.

One has to uncomment

```
///#define LoopGAMMA
```

to force micrOMEGAs to calculate point like gamma signal.

Function `loopGamma(&vcs_gz,&vcs_gg)` calculates annihilation rates `vcs_gz` and `vcs_gg` [cm^3/s]. For example for IDM model with `data1.par`

Gamma ray lines:

$E=5.97\text{E}+02[\text{GeV}]$ $vcs(Z,A)= 1.58\text{E}-28[\text{cm}^3/\text{s}]$, $\text{flux}=4.91\text{E}-14[\text{cm}^2 \text{s}]^{-1}$

$E=6.00\text{E}+02[\text{GeV}]$ $vcs(A,A)= 5.37\text{E}-29[\text{cm}^3/\text{s}]$, $\text{flux}=3.33\text{E}-14[\text{cm}^2 \text{s}]^{-1}$

Calculated cross sections can be compared with experiments for search of single lines in photon spectrum

The `loopGamma` function is not available automatically for models implemented by the user.

Halo profile

DM distribuion is defined by DM density at Sun, parameter `rhoDM` and halo profile. By default micrOMEGAs uses Zhao profile

$$F_{halo}(r) = \left(\frac{R_{\odot}}{r} \right)^{\gamma} \left(\frac{r_c^{\alpha} + R_{\odot}^{\alpha}}{r_c^{\alpha} + r^{\alpha}} \right)^{\frac{\beta-\gamma}{\alpha}}$$

with `alpha=1`, `beta=3` `rc=20kpc`.

`setProfileZhao`(`alpha`,`beta`,`gamma`,`rc`) change these parameters.

`setHaloProfile`(`F`) allows to substitute any profile presented by function `F(r)`

The command `setHaloProfile(hProfileZhao)` sets back the Zhao profile

Photon flux

`gammaFluxTab(fi,dfi,sigmav,Sg,Sobs)`

`fi` is the angle between the line of sight and the center of the galaxy,

`dfi` is half the cone angle which characterizes the detector

resolution (the solid angle is $2\pi(1 - \cos(df\ i))$),

`sigmav` is the annihilation cross section,

`Sg` - photon spectrum at point of annihilation

`Sobs` is tabulated photon flux

`SpectdNdE(E,Sobs)`

gives resulting photon flux in $[1/(\text{GeV cm}^2 \text{s})]$ units

`gammaFlux(fi,dfi,vcs_gz)`

`gammaFlux(fi,dfi,2*vcs_gg)`

return corresponding fluxes for loop induced processes

Antiproton and positron fluxes

- **posiFluxTab**(Emin,sigmav, Se, Sobs)
- **pbarFluxTab**(Emin,sigmav, Sp, Sobs)

The same style as for photons. But depends on propagation parameters

<code>K_dif</code>	0.0112	kpc ² /Myr	The normalized diffusion coefficient
<code>L_dif</code>	4	kpc	Vertical size of the Halo diffu
<code>Delta_dif</code>	0.7		Slope of the diffusion coefficient
<code>Tau_dif</code>	10 ¹⁶	s	Electron energy loss time
<code>Vc_dif</code>	0	km/s	Convective Galactic wind

And finally

solarModulation(Phi, mass, stellarTab, earthTab)

allows to take into account solar modulation effect.

Here **Phi** potential [MeV], **mass** is mass of particle,

stellarTab flux before modulation

earthTab flux after modulation.

[E-print 1004.1092\[hep-ph\]](https://arxiv.org/abs/1004.1092)

Neutrino telescope

micrOMEGAs uses direct detection module to calculate number of DM captured by Sun/Earth.

Captured DM is concentrated in the center of Sun/Earth and neutrino produced in result of DM annihilation can be detected by neutrino telescope experiment ([IceCube](#), [Super-Kamiokande](#), [Baksan](#)).

DM annihilation inside of Sun/Earth is different from annihilation in vacuum. Also there are effects of propagation and oscillation.

For flux of resulting muon neutrinos micrOMEGAs uses tables obtained by [WimpSim](#) package: J. Edso et.al arXiv 0709.3898

Or

[PPPC4DMnu](#): M. Cirelli, et.al. arXiv 1312.6408

Agreement between two sets is not very good.

MicrOMEGAs routine [basicNuSpectra](#) reads these tables depending on

[WIMPSIM](#) flag

[WIMPSIM=1](#) for WimpSim

[WIMPSIM=0](#) for PPC4DMnu

Plots

`displayPlot(title,xName,xMin,xMax,IScale, N, ...)`

displays several curves/histograms on one plot.

`xName` is the name of a variable,

`xMin,xMax` are the lower and upper limits.

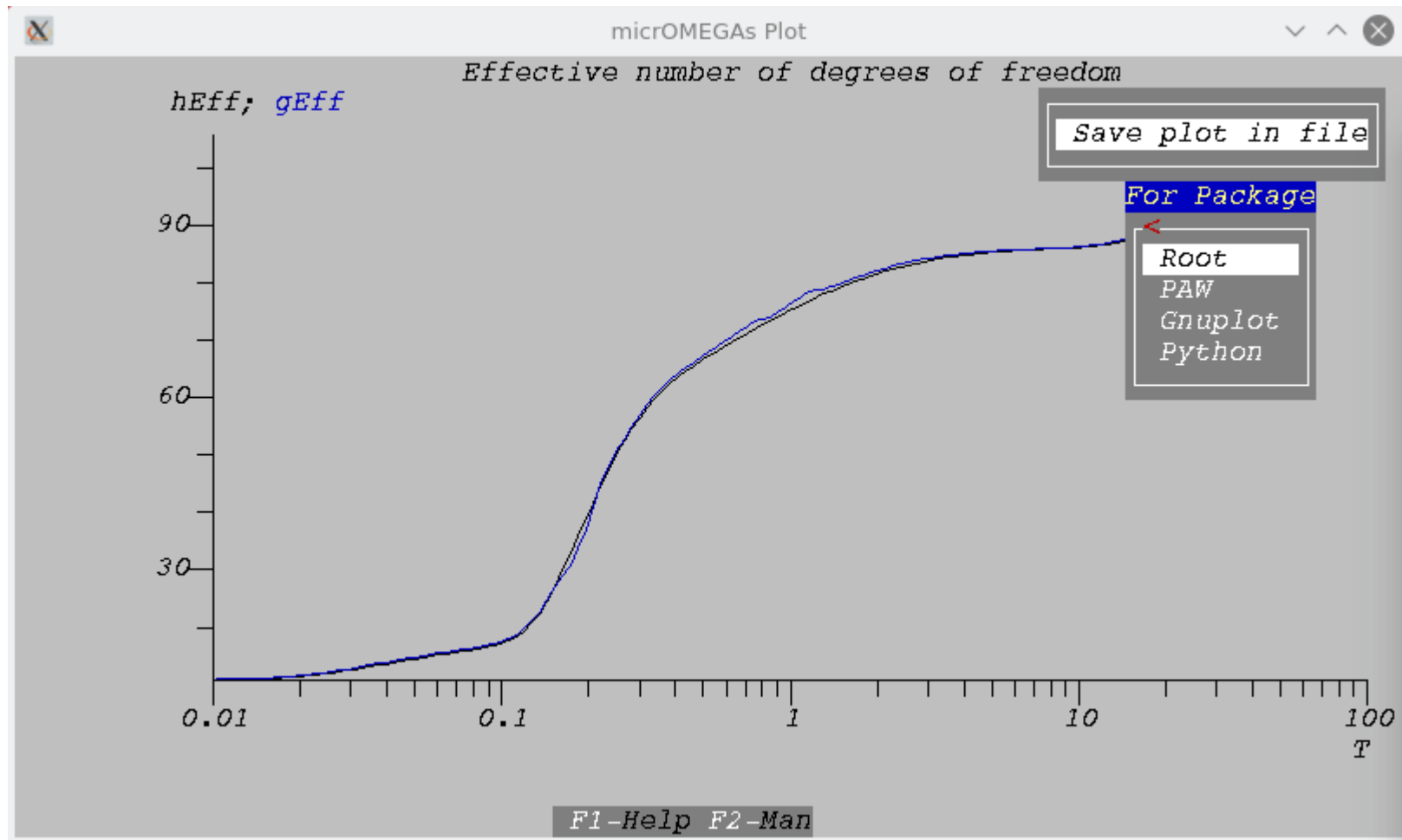
If `IScale` $\neq 0$, a logarithmic scale is used for the x axis.

`N` is the number of curves/histograms to display.

After the parameter `N`, `displayPlot` expects $N \times 4$ parameters, where each tetrad can contain

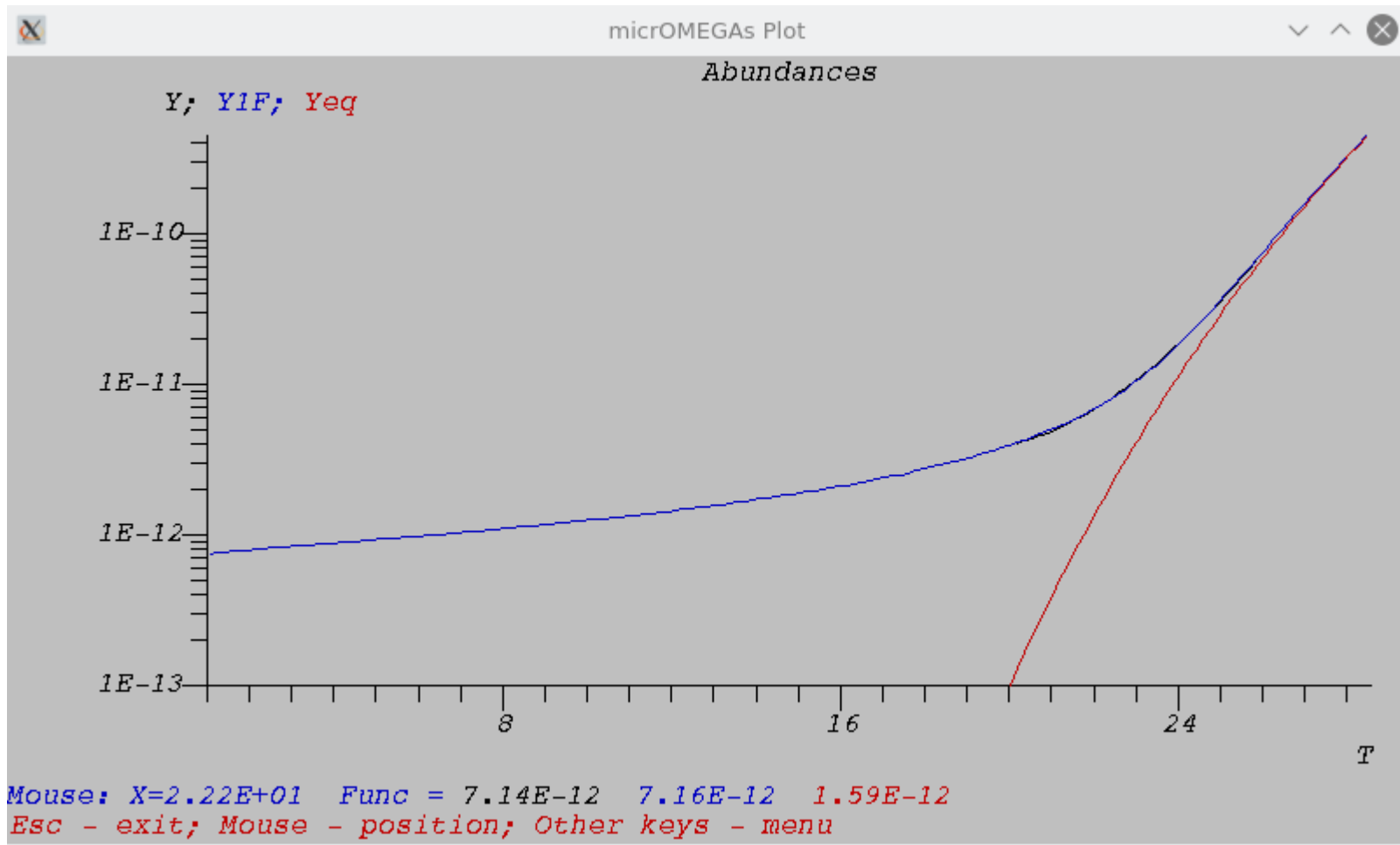
- 1) test label of curve
- 2) dimension of array or 0 for functions
- 3) array of data or a function
- 4) array of error, or address of function arguments, or NULL

Plots: $h_{\text{Eff}}(T)$, $g_{\text{Eff}}(T)$



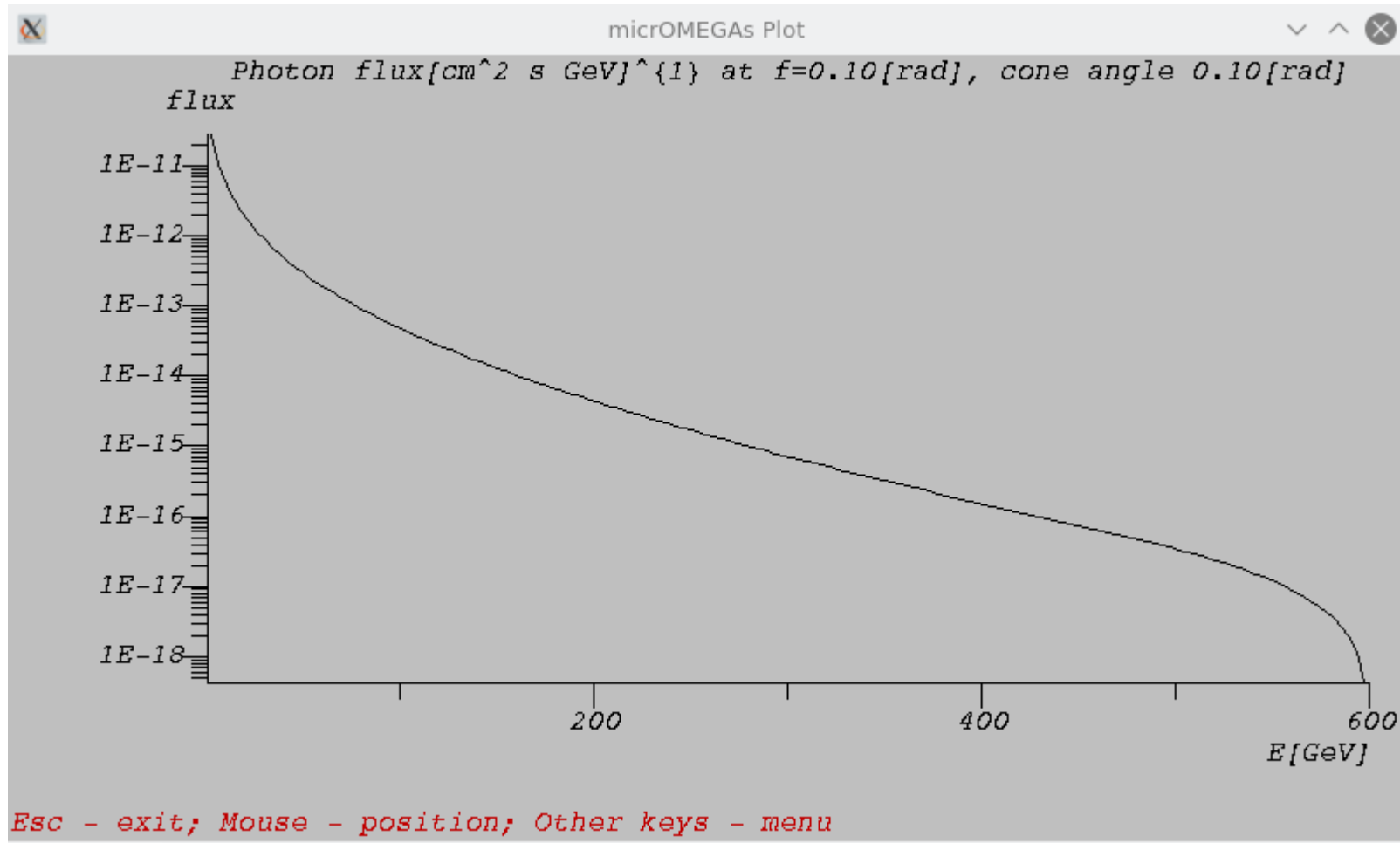
```
displayPlot("Effective number of degrees of freedom","T",0.01,100,  
1,2, "hEff",0,hEff,NULL, "gEff",0,gEff,NULL);
```

Plots: Abundance $Y(T)$, $Y1F(T)$, $Yeq(T)$



```
displayPlot("Abundances","T", 1,Tstart, 0,3,  
"Y",0,YF,NULL, // obtained by darkOmega  
"Y1F",0,Y1F,NULL, // obtained by darkOmega2  
"Yeq",0,Yeq,NULL); // equilibrium abundance
```

Plots: Photon spectrum $\text{SpectdNdE}(E, \text{Flux})$



```
displayPlot(txt,"E[GeV]",Emin,Mcdm,0,1,"flux",0,SpectdNdE,Flux);
```

Here **Flux** is presented by array

External Packages

Directory Packages in intended for external packages

MicrOMEGAs contains several packages which are used for calculation of loop-improved particles spectra in MSSM-like models.

- 1) [SuSpect](#) – for MSSM,
- 2) [NMSSMTools](#) for NMSSM,
- 3) [CpsuperH](#) -for MSSM with complex parameters

Also micrOMEGAs contains [LoopTools](#) used for DM, DM \rightarrow A,A and [Lilith](#) for testing Higgs physics.

Several packages are downloaded automatically when the user calls corresponding functions:

- [HiggsBounds](#) & [HiggsSignals](#) - experimental constraints on Higgs
- [SMODELS](#) – LHC constraints on new particles
- [SPheno](#) & [SoftSUSY](#) - particles spectra of MSSM
- [SuperIso](#) – flavour physics constraints.

SLHAplus[arXiv 1008.0181]: Tools for reading SLHA files

As a rule external packages use SLHA format: hep-ph/0311123, 0801.0045

Routines **slhaRead, slhaVal (read) openAppend, aPrintF (write)**

File with particle spectrum

```
BLOCK MASS # Mass spectrum
# PDG Code mass particle
25 1.15137179E+02 # neutral Higgs
37 1.48428409E+03 # charged Higgs

BLOCK NMIX # Neutralino Mixing Matrix
1 1 9.98499129E-01 # Zn11
1 2 -1.54392008E-02 # Zn12
```

CalcHEP model file

```
slhaRead(file_name, mode)
Mh | slhaVal("MASS",0,1,25)
MHC | slhaVal("MASS",0,1,37)
Zn12 | slhaVal("NMIX",0,2,1,2)
```

Example: SUGRA with SuSpect

```
open | openAppend("suspect2_lha.in")
input1 | aPrintF("Block MODSEL # Select model\n 1 1 # SUGRA\n")
input2 | aPrintF("Block SMINPUTS\n 5 %E#mb(mb)\n 6 %E#mt(pole)\n",MbMb,Mtp)
input3 | aPrintF("BLOCK MINPAR\n 1 %E #m0\n 2 %E #m1/2\n ",Mzero,Mhalf)
input4 | aPrintF("3 %E #tb\n 4 %E #sign(mu)\n 5 %E #A0\n",tb,sgn,A0)
sys | System("$CALCHEP/./Packages/SuSpect_2.41/suspect2.exe")
rd | slhaRead("suspect2_lha.out",0) % mode=4 do not read decays
Mh | slhaVal("MASS", 0 , 1 , 25)
```

SLHAplus updated

Now people use SLHA format more widely than it was proposed by Peter Skands. micrOMEGAs SLHA package was updated correspondingly. For example

Block HiggsBoundsResults

#CHANNELTYPE 1: channel with the highest statistical sensitivity

1	1	328								# channel id
1	2	1								# HBresult
1	3	0.72692779334500290								# obsratio
1	4	1								# ncombined
1	5	(p p)->h+..., h=1 where h is SM-like (CMS-PAS-HIG-12-008)								# channel

slhaSTRFormat("HiggsBoundsResults","1 5 ||%[^]||", **channel**);

Block FOBS # Flavour observables

#	ParentPDG	type	value	q	NDA	ID1	ID2	ID3	...	comment
5	1	2.95061156e-04	0	2	3	22				# BR(b->s gamma)
521	4	8.35442304e-02	0	2	313	22				# Delta0(B->K* gamma)
531	1	3.24270419e-09	0	2	13	-13				# BR(B_s->mu+ mu-)

Bsg= **slhaValFormat**("FOBS", 0., "5 1 %E 0 2 3 22")

Implementation of new models in micrOMEGAs

- The command

`./newProject MODEL`

launched from the root micrOMEGAs directory creates the directory *MODEL*, which contains all files needed to run micrOMEGAs (for example main.c) with the exception of the new model files.

- The new model files in the CalcHEP format should then be included in the sub-directory *MODEL/work/models*. The files needed are

`vars1.mdl, func1.mdl, prtcls1.mdl, lgrng1.mdl extlib1.mdl`

Simple example:

`./newProject IDMcopy`

`cp IDM/work/models/*1.mdl IDMcopy/work/models`

`cp IDM/*.dat IDMcopy`

It should work!

Implementation of new models:

Generation of model files in CalcHEP Format

Model files can be created by mean of

LanHEP , **FeynRules**, **Sarah**

LanHEP [E-print 1412.5016] is included in **micrOMEGAs** package. Each model directory contains **lanhep** subdirectory with source files and with **Makefile** which calls **LanHEP**.

See **LanHEP** manual

[micromegas_X.Y/Packages/LanHEP/manual/man31.pdf](#)

Follow examples presented in any **micrOMEGAs** model.

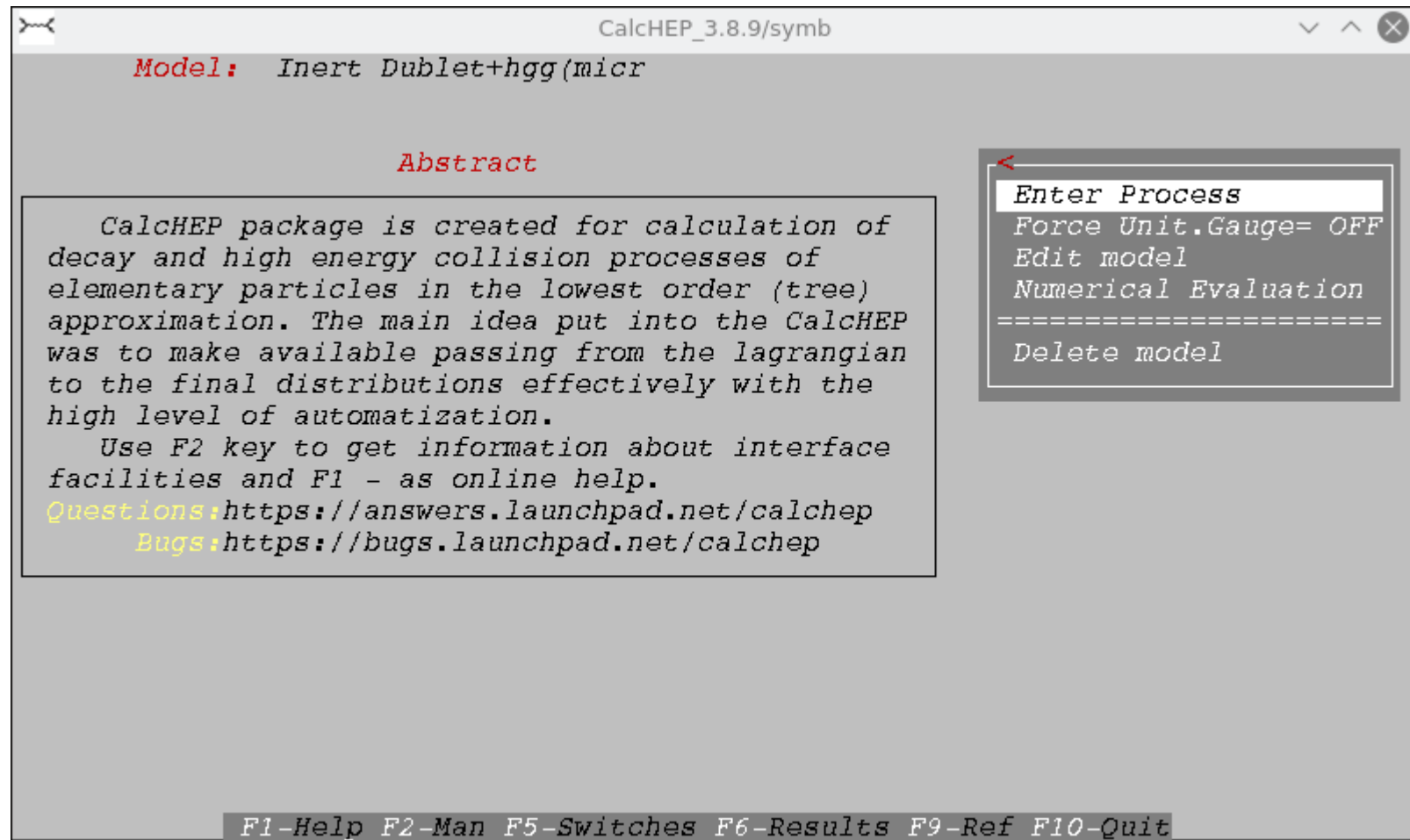
The simplest one is in **IDM/lanhep**

The command **make** launched from the Model directory checks the model and stops with error code if model files do not correspond to CalcHEP requirements.

The user can go to work directory, launch **./calchep** and use menu line

CHECK MODEL

CalcHEP GUI sessions



CalcHEP: Enter process

```
CalcHEP_3.8.9/symb
Model: Inert Doublet+hgg(micr

List of particles (antiparticles)

A(A )- photon          Z(Z )- Z boson        G(G )- gluon
W+(W- )- W boson      ne(Ne )- neutrino    e(E )- electron
nm(Nm )- mu-neutrino  m(M )- muon          nl(Nl )- tau-neutrino
l(L )- tau-lepton     u(U )- u-quark       d(D )- d-quark
c(C )- c-quark        s(S )- s-quark       t(T )- t-quark
b(B )- b-quark        h(h )- Higgs         ~H3      - odd Higgs
~H+(~H-)- Charged Higgs  ~X(~X )- second Higgs

Enter process: ~X,~X -> W+,W-
```

CalcHEP: diagrams

CalcHEP_3.8.9/symb

Delete, On/off, Restore, Latex 1/4

F1 -Help, F2 -Man, PgUp, PgDn, Home, End, # , Esc

CalcHEP: compilation

The image shows two overlapping windows from the CalcHEP 3.8.9 software. The top window, titled 'CalcHEP_3.8.9/symb', displays the model and process settings. The bottom window, titled 'CalcHEP_3.8.9/num', shows the subprocess configuration and a table of results.

CalcHEP_3.8.9/symb

Model: Inert Doublet+hgg(micr

Process: $\sim X, \sim X \rightarrow W^+, W^-$

Feynman diagrams

4 diagrams in 1 subprocesses are constructed.
0 diagrams are deleted.

Squared diagrams

10 diagrams in 1
0 diagrams are d
10 diagrams are c

CalcHEP_3.8.9/num

(sub)Process: $\sim X, \sim X \rightarrow W^+, W^-$
Monte Carlo session: 1

#IT	Cross section[pb]	Error[%]	nCall	Eff.	chi ²
-----	-------------------	----------	-------	------	------------------

CalcHEP_3.8.9/symb menu:

- C code
- C-compiler**
- Edit Linker
- REDUCE code
- MATHEMATICA code

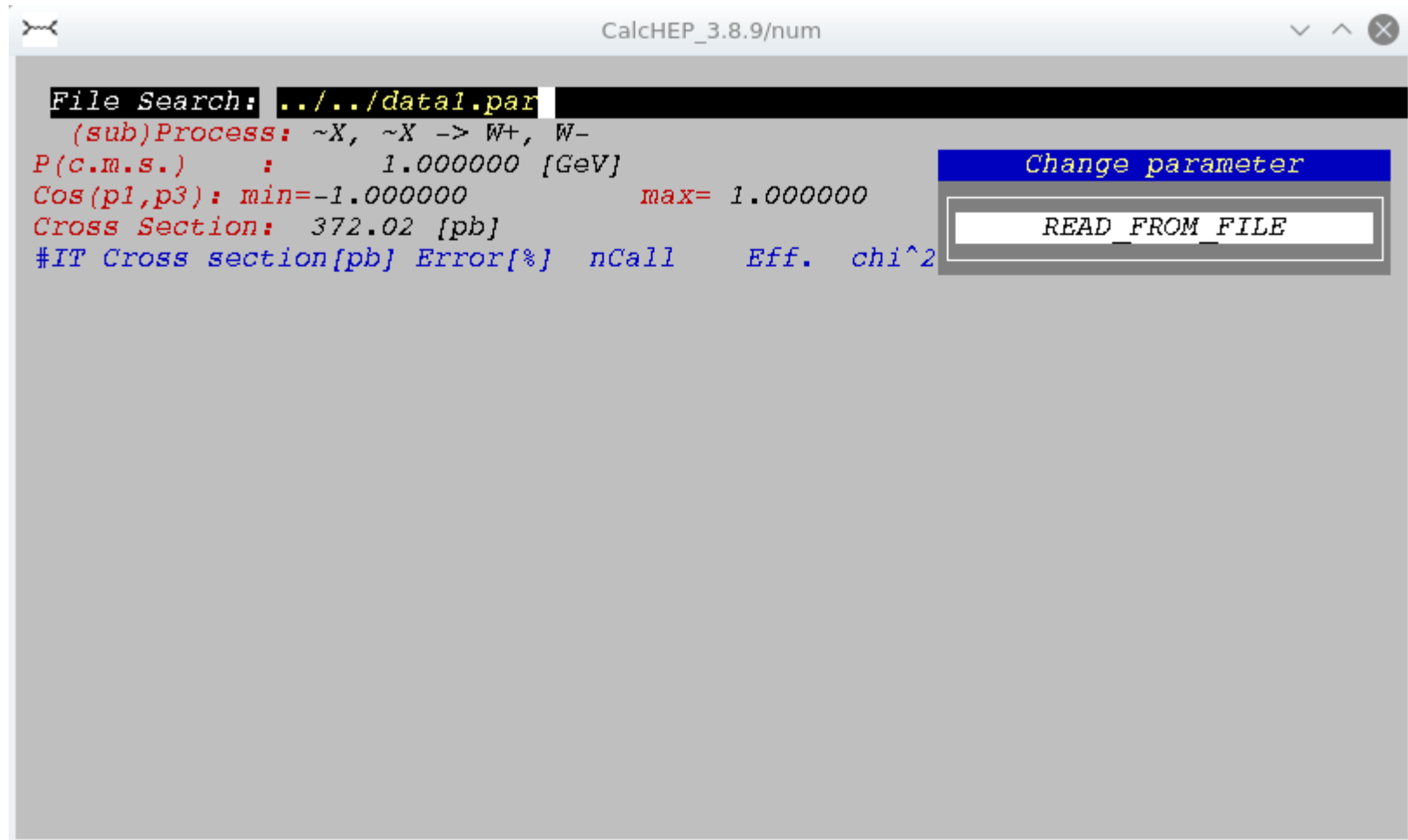
CalcHEP_3.8.9/num menu:

- Subprocess**
- IN state
- Model parameters
- Constraints
- QCD alpha & scales
- Breit-Wigner
- Aliases
- Cuts
- Phase space mapping
- Monte Carlo simulation
- 1D integration

CalcHEP_3.8.9/symb footer: F1-Help F2-Man F3-

CalcHEP_3.8.9/num footer: F1-Help F2-Man F5-Options F6-Results F8-Calc F9-Ref F10-Quit

CalcHEP: parameters



The screenshot shows a window titled "CalcHEP_3.8.9/num". The interface displays the following text:

```
File Search: ../../data1.par  
(sub)Process: ~X, ~X -> W+, W-  
P(c.m.s.) : 1.000000 [GeV]  
Cos(p1,p3): min=-1.000000 max= 1.000000  
Cross Section: 372.02 [pb]  
#IT Cross section[pb] Error[%] nCall Eff. chi^2
```

On the right side of the window, there are two buttons: "Change parameter" (blue background) and "READ FROM FILE" (white background with a black border).

CalcHEP: $v \cdot \sigma$

