

# **CalcHEP version 2.5**

**A.Pukhov**, *SINP MSU, Russia*

**Alexander Belyaev** *Southampton University, GB*

**Neil Christensen** *Michigan State University, US*

**Tools2008**

**Munich, June 30 - July 4**

## **General features of CalcHEP.**

*CalcHEP is a package for **computation** of Feynman diagrams, **integration** over multi-particle phase space, and partonic level event **generation**.*

*Main features are: CalcHEP*

- *Works with generic model;*
- *Calculates symbolically squared Feynman diagrams;*
- *Has user-friendly menu driven interface with on-line help.*
- *Batch mode is based on interactive version (simulation of keyboard hits).*

**WEB disposition:** CalcHEP: <http://theory.sinp.msu.ru/~pukhov/calchep.html>

*Version 2.5 has to appears on WEB in the end of this workshop.*

## **BSM implemented in CalcHEP.**

*During development of micrOMEGAs are realized MSSM(SuSpect, Isajet, SoftSusy, SPheno), NMSSM(NMSSNtools), CPVMSSM(CPSuperH), RHNM*

*The following BSM realizations are known*

*LeptoQuarks(A. Belyaev et al, hep-ph:/0502067)*

*Little Higgs ( A. Belyaev, et al arXiv:hep-ph/0609179)*

*Higgsless 3-site model ( A. Belyaev, S. Chivukula, N. Christensen, E. Simmons et. al. arXiv:hep-ph/arXiv:0708.2588)*

*MUED5, MUED6 ( see Kong WEB page <http://home.fnal.gov/~kckong/mued/>)*

*Almost all these realization were done by means of **LanHEP** package.*

## **CalcHEP development.**

*Last years CalcHEP mainly was developed as a matrix element generator for astro-particle package **micrOMEGAs**.*

*Recently in collaboration with **Alexander Belyaev and Neil Christensen** **CalcHEP** was significantly improved for collider physics applications.*

*Namely it was realized a batch for **multi-processes calculations with decays of outgoing particles and event mixing using parallel calculations on PC-farms**. This job mainly was done by **Alexander Belyaev and Neil Christensen** and will be presented by Neil in the second part of this talk.*

## Other developments ... Models with spin 2 and spin 3/2 particles

*A possibility to work with high spin particles was realized*

Particles								
Full name	P	aP	number	2*spin	mass	width	color	aux
electron	e	E	11	1	0	0	1	
s_electron	~eL	~EL	1000011	0	MSeL	0	1	
Gravitino	~g	~g	1000039	3	MGr	wGr	1	
Massive Graviton	Gh	Gh	2000039	4	MHGr	whGr	1	

*Then interaction can be presented like*

Vertices					Factor	< > Lorentz part
A1	A2	A3	A4	>		
~eL	E	~g		>	$i/\text{Sqrt}2/Mp1$	$ G(m3)*G(p1)$
~EL	~g	e		>	$i/\text{Sqrt}2/Mp1$	$ G(p1)*G(m2)$

## ***Processes with polarized incoming particles.***

*It is realized for incoming massless particles of spin 1/2 and spin 1. One have to use the % symbol before particle name to force CalcHEP to calculate diagrams symbolically using polarized density matrix.*

$e\%, E\% \rightarrow 2*x$

*In numerical session the user can set numerical values for helicities.*

## Automatic width calculation.

*Particle widths included in the model can be calculated automatically using vertices of interaction. It mainly is interesting for BSM particles whose masses unknown.*

Particle									
Full name	P	aP	number	2*spin	mass	width	color	aux	
Light Higgs	h	h	25	0	Mh	!wh	1		
Heavy higgs	H	H	35	0	MHH	!wHh	1		

*The exclamation mark at particle width is a instruction to CalcHEP to generate the code for  $h \rightarrow 2*x$ , and  $H \rightarrow 2*x$  and compile them together with the codes of main processes.*

## Code optimization in CalcHEP

*One problem of CalcHEP is a huge size of codes for 2→4 processes. We have realized a option to reduce significantly size of generated code and accelerate numerical calculation.*

*In general CalcHEP compiles code for diagram which mainly is a polynomial over momenta scalar product with coefficients which are functions on model parameters. Now CalcHEP calculates one time for all diagrams all needed products of scalar products and store them in array. Coefficients at this monomial are calculated one time for session and also stored in some array.*

*So, the code for one diagram is extremely short now.*

```
static const int jumps[30]={266,271,272,274,276,285,293,294,393,395,398
399,409,410,418,419,422,424,429,432,433,435,441,442,459,468,477,478,492
};
for(i=0,R=0 ;i<30;i++){R+=CC[i]*DPmonom[jumps[i]];}
```

*Array jumps is used to bypass zero coefficients.*

## *After MC4BSM3 workshop ... Quadruple Precision*

*We met some physical applications which need calculation with high precision. We have a switch which force high precision calculation.*

*Even standard gcc compiler on PC supports **long double** length of 10.*

*Old platforms like OSF1 work with **long double** length of 16.*

*Modern **Intel** compiler also supports real type **quad** length of 16.*

*CalCHEP\_2.5 supports calculations with high precision real types.*

## Universal SLHA reader.

*The reader for SLHA with arbitrary BLOCK structure is created.*

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

*There is one command to read SLHA file:*

*`slhaRead(char*fileName,int mode)`*

*which downloads in **computer memory** all information and another command which returns the need item*

*`slhaVal(char * BlockName,double Scale, int KeyLength, ...);`*

*There is a possibility to check that the needed item exists: For example*

*`if(slhaValExists("Nmix",2,1,2))Z12=slhaVal("Nmix",100.,2,1,2)); else ...`*

*If there are several blocks with the same name but different scales, `slhaVal` **interpolates** numbers for **Scale** point.*

## **Universal SLHA reader.**

### **Work with *DECAY* items**

```
DECAY 1000037 3.895221E+00 # second chargino(~2+) width
# Branching Nout ID_1 ID_2 ....
1.8877E-03 2 24 1000022 # Br(~2+ -> W+ ~o1)
2.4367E-01 2 25 1000024 # Br(~2+ -> h ~1+)
.....
```

***is organized in the same manner:***

```
int slhaDecayExists(int pNum);
double slhaWidth(int pNum);
double slhaBranch(int pNum,int N, int * nCh);
```

***The is code disposed in CalcHEP/c\_sources/model\_aux/SLHAread.c and can be used out of CalcHEP package.***

***Even Fortran interface is written.***

## Interface with Showering and Hadronization event Generators

*This interface is based on Les Houches Accord for event generators hep-ph/0109068—.*

*But instead of FORTRAN COMMON BLOCKs the information is passed into **text file** according to the later conventions.*

*In result of work of CalcHEP event generator the event\_#.txt files are created in **different directories**.*

```
SCANDIR('dirname')
```

*command opens all event\_#.txt files and reads their headers. It can be called **several times for different directories** for both **scattering and decay events**.*

*After that any **call** of UPEVNT will choose a subprocess according to its **cross section**, read an event record and substitute **recursively all decays** generated.*

***QCD color flows** generated in collisions are continued in **decays tails**.*

## Interface with .... Qnumbers

The main idea of this approach was to share **New Physics/SM physics** between **CalcHEP/PYTHIA**. In principle only **SM particles** should be passed to **PYTHIA**.

From the other side in the event record we keep information about all intermediate particles, in particular their **life time** is saved for subsequent reconstruction of **space-time picture** of event.

So, we have to pass to **PYTHIA** information about quantum numbers of **non-SM particles**.

It is done via **QNUMBERS** (P. Skands et al, hep-ph/0311123) block generated by **CalcHEP** automatically and include description of all non-SM particles:

```
BLOCK QNUMBERS 1000045 # ~chi_50
 1  0  # 3 times electric charge
 2  2  # number of spin states (2S+1)
 3  1  # color rep (1: singlet, 3: triplet, 8: octet)
 4  0  # Particle/Antiparticle distinction (0=own anti)
```

## **The Batch job**

*The preliminary calls of CalcHEP for cross section calculation and one channels event generations and subsequent event mixing with recursive decays can be realized by one **BATCH** command which will be presented by **Neil Christensen** now.*