# **SPheno**, a program for calculating supersymmetric spectra, SUSY particle decays and SUSY particle production at $e^+e^-$ colliders

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#### Abstract

SPheno is a program that accurately calculates the supersymmetric particle spectrum within a high scale theory, such as minimal supergravity, gauge mediated supersymmetry breaking, anomaly mediated supersymmetry breaking, or string effective field theories. An interface exists for an easy implementation of other high scale models. The program solves the renormalization group equations numerically to two–loop order with user-specified boundary conditions. The complete one–loop formulas for the masses are used which are supplemented by two–loop contributions in case of the neutral Higgs bosons and the  $\mu$  parameter. The obtained masses and mixing matrices are used to calculate decay widths and branching ratios of supersymmetric particles as well as of Higgs bosons,  $b \to s \gamma$ ,  $\Delta \rho$  and  $(g-2)_{\mu}$ . Moreover, the production cross sections of all supersymmetric particle as well as Higgs bosons at  $e^+e^-$  colliders can be calculated including initial state radiation and longitudinal polarization of the incoming electrons/positrons. The program is structured such that it can easily be extend to include non-minimal models and/or complex parameters. Starting with version 2.2.2 the SLHA convention as well as the SPA convetion are supported.

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# 1 Introduction

Supersymmetry (SUSY) [1, 2, 3] provides an attractive extension of the Standard Model (SM). It provides a qualitative understanding of various phenomena in particle physics: It stabilizes the gap between the Grand Unification scale / Planck scale and the electroweak

scale [4]. It allows the unification of the three gauge couplings at a scale  $M_U \simeq 2 \cdot 10^{16}$  GeV in a straight forward way [5]. The large top mass generates radiative electroweak symmetry breaking [6]. In addition it provides the lightest supersymmetric particle as a cold dark matter candidate [7]. Therefore, the search for supersymmetric particles is one of the main topics in the experimental program of present and future high energy colliders [8, 9, 10, 11].

The Minimal Supersymmetric Standard Model (MSSM) consists of taking the Standard Model and adding the corresponding supersymmetric partners [3]. In addition a second Higgs doublet is needed to obtain an anomaly-free theory. The second Higgs doublet is also needed to give mass to u-type quarks and down-type quarks at the same time. The MSSM in its most general from contains more than 100 unknown parameters [12] which are clearly to many for an exhaustive study. This number drastically reduces if one embeds the MSSM in a high scale theory, such as minimal supergravity theories [13], gauge mediated supersymmetry breaking[14], or anomaly mediated supersymmetry breaking[15]. There is not yet a theoretical preferred scheme for supersymmetry breaking. For this reason it is important to know whether the precision of on-going and future experiments is sufficient: (i) To distinguish between the various schemes. (ii) To which extend it is possible to reconstruct the underlying theory. It has been demonstrated that the expected experimental accuracies at future  $e^+e^-$  colliders complemented with data from the LHC allow for a successful reconstruct of such an underlying supersymmetric high scale theory [16]. Connected with these questions is the question if the theoretical accuracy matches the experimental one. The present version of the program SPheno<sup>1</sup> is thought as a further step in getting accurate theoretical results to match finally the experimental precision.

In the view of ongoing and future experiments it is highly desirable to have various and independent tools at hand performing the calculation of the supersymmetric spectrum, of decay widths, of branching ratios and of production cross sections. This allows for a cross check of the tools and by comparing the implemented methods and approximations one can also get a rough understanding of the theoretical error. SPheno is a program performing an accurate calculation of the supersymmetric spectrum, of the branching ratios of supersymmetric particles and the Higgs bosons and of the production cross sections of supersymmetric particles and the Higgs bosons in  $e^+e^-$  annihilation including longitudinal beam polarization. Moreover, the spectrum is used to calculate the branching of the rare decay  $b \to s \gamma$ , the supersymmetric contributions to the anomalous magnetic of the muon  $a_{\mu}$  as well as supersymmetric contributions to the  $\rho$  parameter.

For the calculation of the spectrum the programs ISAJET [17], SOFTSUSY [18] and SUSPECT [19] are widely used. A comparison of the results among these programs and with SPheno is given in [20]. The calculation of the branching ratios of supersymmetric particles as well as the production cross sections in  $e^+e^-$  annihilation can be done with SPheno, ISAJET [17], SPYTHIA [21], SUSYGEN [22] and SDECAY [23]. A comparison of the results of these programs will be given in a future paper.

SPheno has been written in Fortran 90. The main focus has been on accuracy and on stable numerical results and less on speed. However, on a modern PC a typical running time is in the order of one second. The calculation is done using two-loop renormalization group equations (RGEs) [24], complete one-loop correction to all SUSY and Higgs masses [25] supplemented by the 2-loop corrections to the neutral Higgs bosons [26, 27] and to the  $\mu$  parameter [27]. The present version of SPheno does all calculations for real parameters

<sup>&</sup>lt;sup>1</sup>SPheno stands for **S**upersymmetric **Pheno**menology

neglecting the flavour structure in the fermion as well as in the sfermion sector. Decay widths and cross sections are calculated using tree-level formulas. However, the couplings involved are running couplings and thus important numerical effects of higher order corrections are already taken into account. The program has been structured in such a way that the future inclusion of complex phases and mixing between the generations has already been considered in the design of the interfaces as well as in the definition of the various variables. Moreover, extensions of the MSSM, e.g. models with broken R-parity, can be implemented easily.

The aim of this paper is to provide a manual of the program, version 2.0, to describe the approximations used and to display the results of a run. In Sect. 2 we will summarize the MSSM parameters and we give the tree-level formulas for the supersymmetric particles. Moreover, a short summary of the implemented high scale models is given. In Sect. 3 we list the implemented decay modes of supersymmetric particles and the Higgs bosons. We also discuss shortly the approximations used. In Sect. 4 we present the implemented cross sections in  $e^+e^-$  annihilation. In Sect. 5 we discuss give details on the implemented low energy constraints. In Sect. 6 we discuss the implemented algorithm in some detail. In Sect. 7 the main program is presented in detail providing the necessary information so that this program can be easily adapted to the user's requirement. In the appendices we discuss the possible switches for influencing the program as well as a detailed discussion of possible input files. This can be done using SPheno specific files or by using the SUSY Les Houches Accord (SLHA) [28]. Moreover, we list the output of the program for a typical example. The source code as well as precompiled a version of the program can be obtained from the author via email: porod@physik.unizh.ch or porod@ific.uv.es; or it can be downloaded from http://www-theorie.physik.unizh.ch/~porod/SPheno.html.

## 2 MSSM parameters, particle spectrum, and Models

In this section we fix our notation concerning the parameters and present the tree-level formulas for the masses as well as the mixing matrices. In the following we assume that the physical masses are ordered:  $m_i \leq m_j$  if i < j except for the sfermions as explained below. We also give a short overview over various high scale models which are implemented in the program.

## 2.1 Ingredients for the Lagrangian

The pure supersymmetric Lagrangian is specified by the Kähler potential giving the gauge interactions and by the Superpotential W giving the Yukawa interactions:

$$W = \epsilon_{ab} \left( Y_{ij}^L \hat{L}_i^a \hat{H}_1^b \hat{E}_j^c + Y_{ij}^D \hat{Q}_i^a \hat{H}_1^b \hat{D}_j^c + Y_{ij}^U \hat{Q}_i^b \hat{H}_2^a \hat{U}_j^c - \mu \hat{H}_1^a \hat{H}_2^b \right)$$
(1)

where  $\hat{L}$ ,  $\hat{E}$ ,  $\hat{Q}$ ,  $\hat{D}$ , and  $\hat{U}$  denote the matter superfields. The  $SU(2)_L$  representation indices are denoted by a, b = 1, 2 and the generation indices by i, j = 1, 2, 3;  $\epsilon_{ab}$  is the totally antisymmetric tensor with  $\epsilon_{12} = 1$ . Note that the sign of  $\mu$  is identical to the one in ISAJET [17] and SOFTSUSY [18] but opposite to the convention in [25]. Presently, real Yukawas  $Y^L$ ,  $Y^D$ ,  $Y^U$  only are included. They and the gauge couplings  $g_i$  are  $\overline{DR}$  quantities.  $g_1$  is defined in the Grand Unification normalization  $g_1 = \sqrt{5/3}g'$  where g' is the Standard Model hypercharge gauge coupling.

The next ingredient is the soft SUSY breaking Lagrangian, which is given by mass terms for the gauginos

$$L_{soft,1} = \frac{1}{2} \left( M_1 \tilde{B} \tilde{B} + M_2 \tilde{W}_a \tilde{W}^a + M_3 \tilde{g}_\alpha \tilde{g}^\alpha \right) + h.c. , \qquad (2)$$

mass terms for scalar matter fields and Higgs fields

$$L_{soft,2} = -M_{H_1}^2 H_{1a}^* H_1^a - M_{H_2}^2 H_{2a}^* H_2^a - M_{\tilde{L},ij}^2 \tilde{L}_{ia}^* \tilde{L}_j^a - M_{\tilde{E},ij}^2 \tilde{E}_i^* \tilde{E}_j -M_{\tilde{Q},ij}^2 \tilde{Q}_{ia}^* \tilde{Q}_j^a - M_{\tilde{U},ij}^2 \tilde{U}_i^* \tilde{U}_j - M_{\tilde{D},ij}^2 \tilde{D}_i^* \tilde{D}_j$$
(3)

and trilinear couplings of scalar matter fields and Higgs fields

$$L_{soft,3} = -\epsilon_{ab} \left( A_{ij}^L \tilde{L}_i^a H_1^b \tilde{E}_j^* + A_{ij}^D \tilde{Q}_i^a H_1^b \tilde{D}_j^* + A_{ij}^U \tilde{Q}_i^b H_2^a \tilde{U}_j^* - B\mu H_1^a H_2^b \right) + h.c.$$
(4)

## 2.2 Masses and Mixing Matrices

The masses of the various particles are induced by the soft SUSY breaking parameters and the vacuum expectation values  $v_i$  of the neutral Higgs fields  $v_i = \langle H_i^0 \rangle$ . The ratio of the vacuum expectation values is denoted by  $\tan \beta = v_2/v_1$ . The sum of the vacuum expectation values (vevs) squared is fixed by the gauge boson masses:

$$m_W^2 = \frac{1}{4}g^2(v_1^2 + v_2^2), \qquad m_Z^2 = \frac{1}{4}(g^2 + g'^2)(v_1^2 + v_2^2)$$
 (5)

Neglecting the mixing between different generations, the Standard Model fermion masses are given by:

$$m_{u_i} = \frac{1}{\sqrt{2}} Y_{ii}^U v_2, \qquad m_{d_i} = \frac{1}{\sqrt{2}} Y_{ii}^D v_1, \qquad m_{l_i} = \frac{1}{\sqrt{2}} Y_{ii}^L v_1$$
 (6)

for u-quarks, d-quarks and leptons, respectively.

The gluino mass is given by  $m_{\tilde{g}} = |M_3|$ . The charginos are combination of the charged winos  $\tilde{w}^{\pm} = (\tilde{w}^1 \mp i\tilde{w}^2)/\sqrt{2}$  and the charged higgsinos  $\tilde{h}_1^-, \tilde{h}_2^+$ . The Lagrangian contains the chargino mass term  $-(\tilde{\psi}^-)^T X \tilde{\psi}^+$  where  $\psi^- = (-i\tilde{w}^-, \tilde{h}_1^-)^T, \ \psi^+ = (-i\tilde{w}^+, \tilde{h}_2^+)^T$  and

$$X = \begin{pmatrix} M & \frac{g}{\sqrt{2}}v_2 \\ \frac{g}{\sqrt{2}}v_1 & \mu \end{pmatrix}. \tag{7}$$

The matrix is diagonalized by two unitary matrices U and V:

$$M_{D,\tilde{\chi}^{\pm}} = U^* X V^{-1} \,.$$
 (8)

The neutral gauginos  $\tilde{b}$ ,  $\tilde{w}^3$  as well as the neutral higgsinos  $\tilde{h}^0_1$ ,  $\tilde{h}^0_2$  form the neutralinos. In the basis  $\tilde{\psi}^0 = (i\tilde{b}, i\tilde{w}^3, \tilde{h}^0_1, \tilde{h}^0_2)^T$  one finds the mass term  $-(\tilde{\psi}^0)^T Y \tilde{\psi}^0$  with

$$Y = \begin{pmatrix} M_1 & 0 & -\frac{g'}{2}v_1 & \frac{g'}{2}v_2 \\ 0 & M_2 & \frac{g}{2}v_1 & -\frac{g}{2}v_2 \\ -\frac{g'}{2}v_1 & \frac{g}{2}v_1 & 0 & -\mu \\ \frac{g'}{2}v_2 & -\frac{g}{2}v_2 & -\mu & 0 \end{pmatrix}$$
(9)

This matrix is diagonalized by an unitary matrix N:

$$M_{D,\tilde{\chi}^0} = N^* Y N^{\dagger} \,. \tag{10}$$

The CP-even electroweak eigenstates  $(H_1^0, H_2^0)$  are rotated by the angle  $\alpha$  into the Higgs mass eigenstates  $(h^0, H^0)$  as follows:

$$\begin{pmatrix} h^0 \\ H^0 \end{pmatrix} = \begin{pmatrix} -\sin\alpha & \cos\alpha \\ \cos\alpha & \sin\alpha \end{pmatrix} \begin{pmatrix} H_1^0 \\ H_2^0 \end{pmatrix}$$
 (11)

with  $m_{h^0} < m_{H^0}$ . The CP-odd and the charged Higgs masses are given by

$$m_{A^0}^2 = B \mu (\tan \beta + \cot \beta), \qquad m_{H^+}^2 = m_{A^0}^2 + m_W^2$$
 (12)

at tree level.

Neglecting generation mixing, the sneutrino masses are given by:

$$M_{\tilde{\nu}_i}^2 = M_{\tilde{L}_i i}^2 + \frac{1}{2} m_Z^2 \cos 2\beta \tag{13}$$

The other sfermion mass matrices are  $2 \times 2$  matrices:

$$M_{\tilde{l},i}^{2} = \begin{pmatrix} M_{\tilde{L},ii}^{2} - \left(\frac{1}{2} - s_{W}^{2}\right) c_{2\beta} m_{Z}^{2} + m_{l,i}^{2} & \frac{1}{\sqrt{2}} \left(v_{1} (A_{ii}^{L})^{*} - \mu Y_{ii}^{L} v_{2}\right) \\ \frac{1}{\sqrt{2}} \left(v_{1} A_{ii}^{L} - (\mu Y_{ii}^{L})^{*} v_{2}\right) & M_{\tilde{E},ii}^{2} - s_{W}^{2} c_{2\beta} m_{Z}^{2} + m_{l,i}^{2} \end{pmatrix}$$

$$M_{\tilde{u}}^{2} = \begin{pmatrix} M_{\tilde{Q},ii}^{2} + \left(\frac{1}{2} - \frac{2}{3} s_{W}^{2}\right) c_{2\beta} m_{Z}^{2} + m_{u,i}^{2} & \frac{1}{\sqrt{2}} \left(v_{2} (A_{ii}^{U})^{*} - \mu Y_{ii}^{U} v_{1}\right) \\ \frac{1}{\sqrt{2}} \left(v_{2} A_{ii}^{U} - (\mu Y_{ii}^{U})^{*} v_{1}\right) & M_{\tilde{U},ii}^{2} + \frac{2}{3} s_{W}^{2} c_{2\beta} m_{Z}^{2} + m_{u,i}^{2} \end{pmatrix}$$

$$M_{\tilde{d}}^{2} = \begin{pmatrix} M_{\tilde{Q},ii}^{2} - \left(\frac{1}{2} - \frac{1}{3} s_{W}^{2}\right) c_{2\beta} m_{Z}^{2} + m_{d,i}^{2} & \frac{1}{\sqrt{2}} \left(v_{1} (A_{ii}^{D})^{*} - \mu Y_{ii}^{D} v_{2}\right) \\ \frac{1}{\sqrt{2}} \left(v_{1} A_{ii}^{D} - (\mu Y_{ii}^{D})^{*} v_{2}\right) & M_{\tilde{D},ii}^{2} - \frac{1}{3} s_{W}^{2} c_{2\beta} m_{Z}^{2} + m_{d,i}^{2} \end{pmatrix}$$

$$(14)$$

$$M_{\tilde{u}}^{2} = \begin{pmatrix} M_{\tilde{Q},ii}^{2} + \left(\frac{1}{2} - \frac{2}{3}s_{W}^{2}\right)c_{2\beta}m_{Z}^{2} + m_{u,i}^{2} & \frac{1}{\sqrt{2}}\left(v_{2}(A_{ii}^{U})^{*} - \mu Y_{ii}^{U}v_{1}\right) \\ \frac{1}{\sqrt{2}}\left(v_{2}A_{ii}^{U} - (\mu Y_{ii}^{U})^{*}v_{1}\right) & M_{\tilde{U},ii}^{2} + \frac{2}{3}s_{W}^{2}c_{2\beta}m_{Z}^{2} + m_{u,i}^{2} \end{pmatrix}$$
(15)

$$M_{\tilde{d}}^{2} = \begin{pmatrix} M_{\tilde{Q},ii}^{2} - \left(\frac{1}{2} - \frac{1}{3}s_{W}^{2}\right)c_{2\beta}m_{Z}^{2} + m_{d,i}^{2} & \frac{1}{\sqrt{2}}\left(v_{1}(A_{ii}^{D})^{*} - \mu Y_{ii}^{D}v_{2}\right) \\ \frac{1}{\sqrt{2}}\left(v_{1}A_{ii}^{D} - (\mu Y_{ii}^{D})^{*}v_{2}\right) & M_{\tilde{D},ii}^{2} - \frac{1}{3}s_{W}^{2}c_{2\beta}m_{Z}^{2} + m_{d,i}^{2} \end{pmatrix}$$
(16)

where  $c_{2\beta} = \cos 2\beta$  and  $s_W^2 = \sin^2 \theta_W$ . These matrices are diagonalized by  $2 \times 2$  matrices  $R_{\tilde{f},i}$  with

$$m_{\tilde{f}}^2 = R_{\tilde{f}} M_{\tilde{f}}^2 R_{\tilde{f}}^{\dagger} \tag{17}$$

Sfermions are first ordered according to the generation and inside a generation according to their masses. For example, in the slepton sector the ordering is  $\tilde{e}_1$ ,  $\tilde{e}_2$ ,  $\tilde{\mu}_1$ ,  $\tilde{\mu}_2$ ,  $\tilde{\tau}_1$ ,  $\tilde{\tau}_2$  and similarly for squarks.

#### 2.3 High scale models

In this section we summarize the key ingredients of the high scale models implemented in SPheno. We also present the formulas for the boundary conditions in the various models. In all cases the modulus  $|\mu|$  is determined by requiring correct radiative symmetry breaking. At tree level the corresponding formula reads as:

$$|\mu|^2 = \frac{1}{2} \left[ \tan 2\beta \left( M_{H_2}^2 \tan \beta - M_{H_1}^2 \cot \beta \right) - m_Z^2 \right]. \tag{18}$$

Moreover, in all cases the high scale parameters are supplemented by the sign of  $\mu$  and  $\tan \beta$ .

#### 2.3.1Minimal Supergravity

The minimal supergravity (mSUGRA) scenario is characterized by a set of universal parameters [13, 2] at the GUT scale  $M_{GUT}$ : the gaugino mass parameter  $M_{1/2}$ , the scalar mass parameter  $M_0$ , and the trilinear coupling  $A_0$ :

$$M_i(M_{GUT}) = M_{1/2} \tag{19}$$

$$M_{\tilde{j}}^2(M_{GUT}) = M_0^2$$
 (20)  
 $A_i(M_{GUT}) = A_0 Y_i(M_{GUT})$  (21)

$$A_i(M_{GUT}) = A_0 Y_i(M_{GUT}) (21)$$

#### Minimal Supergravity including right handed neutrinos 2.3.2

In addition to the parameters of the mSUGRA model above the following parameters appear in this case:  $m_{\nu_R}$ , a common mass for all right handed neutrinos, and  $m_{\nu_i}$  (i=1,2,3), the light neutrino masses. In this case the MSSM RGEs are run up to the scale  $m_{\nu_R}$  where the neutrino Yukawa couplings are calculated using the formula  $Y_{\nu,i} = \sqrt{m_{\nu_R} m_{\nu_i}}/v_2$ . In the range between  $m_{\nu_R}$  and  $M_{GUT}$  the effect of neutrino Yukawa couplings is included in the RGEs of gauge and Yukawa couplings. At the GUT-scale the right sneutrino mass parameters as well as the trilinear coupling  $A_{\nu,i}$  are given by:

$$M_{\tilde{R}}^2(M_{GUT}) = M_0^2 (22)$$

$$A_{\nu,i}(M_{GUT}) = A_0 Y_{\nu,i}(M_{GUT}) \tag{23}$$

The corresponding RGEs are used in the running from  $M_{GUT}$  to  $m_{\nu_R}$ . At the scale  $m_{\nu_R}$  the neutrino Yukawa couplings  $Y_{\nu,i}$ , the trilinear couplings  $A_{\nu,i}$  and the soft masses  $M_{R,i}^2$  for the right sneutrinos are taken out of the RGEs and below the  $m_{\nu_R}$  the usual set of MSSM RGEs are used.

#### Gauge Mediated Supersymmetry Breaking 2.3.3

Gauge mediated supersymmetry breaking [14, 29] (GMSB) is characterized by the mass  $M_M \sim \langle S \rangle$  of the messenger fields and the mass scale  $\Lambda = \langle F_S \rangle / \langle S \rangle$  setting the size of the gaugino and scalar masses. The gaugino masses

$$M_i(M_M) = (N_5 + 3N_{10})g(\Lambda/M_M)\alpha_i(M_M)\Lambda$$
(24)

are generated by loops of scalar and fermionic messenger component fields;  $N_i$  is the multiplicity of messengers in the  $5 + \overline{5}$  and  $10 + \overline{10}$  vector-like multiplets, and

$$g(x) = \frac{1+x}{x^2}\log(1+x) + (x \to -x)$$
 (25)

is the messenger-scale threshold function [30] which approaches unity for  $\Lambda \ll M_M$ . Masses of the scalar fields in the visible sector are generated by 2-loop effects of gauge/gaugino and messenger fields:

$$M_{\tilde{j}}^{2}(M_{M}) = 2(N_{5} + 3N_{10})f(\Lambda/M_{M})\sum_{i=1}^{3}k_{i}C_{j}^{i}\alpha_{i}^{2}(M_{M})\Lambda^{2}$$
(26)

with  $k_i = 1, 1, 3/5$  for SU(3), SU(2), and U(1), respectively; the coefficients  $C_i^i$  are the quadratic Casimir invariants, being 4/3, 3/4, and  $Y^2/4$  for the fundamental representations  $\tilde{j}$ 

in the groups i = SU(3), SU(2) and U(1), with  $Y = 2(Q-I_3)$  denoting the usual hypercharge; also the threshold function [30]

$$f(x) = \frac{1+x}{x^2} \left[ \log(1+x) - 2\text{Li}_2\left(\frac{x}{1+x}\right) + \frac{1}{2}\text{Li}_2\left(\frac{2x}{1+x}\right) \right] + (x \to -x)$$
(27)

approaches unity for  $\Lambda \ll M_M$ . As evident from Eq. (26) scalar particles with identical Standard–Model charges squared have equal masses at the messenger scale  $M_M$ . In the minimal version of GMSB, the A parameters are generated at 3-loop level and they are practically zero at  $M_M$ . However, the program permits to set a value for  $A_0$  different from zero but universal for all sfermions.

## 2.3.4 Anomaly Mediated Supersymmetry Breaking

In anomaly mediated supersymmetry breaking (AMSB) the SUSY breaking is transmitted from the hidden sector to the visible sector via the super–Weyl anomaly [15]. The soft SUSY breaking parameters are given by:

$$M_a = \frac{1}{q_a} \beta_a m_{3/2} \tag{28}$$

$$A_i = \beta_{Y_i} m_{3/2} \tag{29}$$

$$M_j^2 = \frac{1}{2} \dot{\gamma}_j m_{3/2}^2 \tag{30}$$

where  $\beta_a$  and  $\beta_{Y_i}$  are the beta functions of gauge and Yukawa couplings, respectively. The  $\gamma_j$  are the anomalous dimensions of the corresponding matter superfield and  $m_{3/2}$  is the gravitino mass. Equation (30) leads to negative mass squared for the sleptons which is phenomenologically not acceptable. There are several possibilities to solve this problem [31] and the simplest one is to add a common scalar mass  $M_0$  so that eq. (30) reads as

$$M_j^2 = M_0^2 + \frac{1}{2}\dot{\gamma}_j m_{3/2}^2 (31)$$

This extension has been implemented in the program.

#### 2.3.5 String Effective Field Theories

Four-dimensional strings naturally give rise to a minimal set of fields for inducing supersymmetry breaking; they play the rôle of the fields in the hidden sectors: the dilaton S and the moduli  $T_m$  chiral superfields which are generically present in large classes of 4-dimensional heterotic string theories. The vacuum expectation values of S and  $T_m$ , generated by genuinely non-perturbative effects, determine the soft supersymmetry breaking parameters [32, 33].

In the following we assume that all moduli fields get the same vacuum expectation values and that they couple in the same way to matter fields. Therefore, we omit the index m and take only one moduli field T. The properties of the supersymmetric theories are quite different for dilaton and moduli dominated scenarios as discussed in [32, 33]. The mass scale of the supersymmetry parameters is set by the gravitino mass  $m_{3/2}$ .

In the program we implemented the complete 1-loop formulas given in [33]. Three classes of models are implemented in the program: two versions of OII compactification defined by

the sets A and B of boundary conditions in [33] as well as an OI compactification scheme. For the implementation of the OI compactification scheme we have used formulas Eqs. (3.21) – (3.23) of [33]:

$$M_{i} = -g_{i}^{2} m_{3/2} \left\{ \sqrt{3} \sin \theta + \left[ b_{i} + s \sqrt{3} \sin \theta g_{s}^{2} \left( C_{i} - \sum_{j} C_{i}^{j} \right) + 2 t \cos \theta G_{2}(t) \left( \delta_{GS} + b_{i} - 2 \sum_{j} C_{i}^{j} (1 + n_{j}) \right) \right] / 16 \pi^{2} \right\}$$

$$M_{j}^{2} = m_{3/2}^{2} \left\{ \left( 1 + n_{j} \cos^{2} \theta \right) + 2 \sqrt{3} s \sin \theta \left[ \sum_{i} \gamma_{j}^{i} g_{i}^{2} - \frac{1}{2 s} \sum_{km} \gamma_{j}^{km} \right] + \gamma_{j} + 2 t \cos \theta G_{2}(t) \sum_{km} \gamma_{j}^{km} (n_{j} + n_{k} + n_{m} + 3) \right\}$$

$$A_{jkm} = m_{3/2} \left[ -\sqrt{3} \sin \theta - 2 t \cos \theta (n_{j} + n_{k} + n_{m} + 3) G_{2}(t) + \gamma_{j} + \gamma_{k} + \gamma_{m} \right]$$

$$(34)$$

 $s = \langle S \rangle$  is the vacuum expectation values of the dilaton field.  $t = \langle T \rangle / m_{3/2}$  is the vacuum expectation value of the moduli field(s), and  $G_2(t) = 2\zeta(t) + 1/2t$  is the non-holomorphic Eisenstein function with  $\zeta$  denoting the Riemann zeta function.  $\delta_{GS}$  is the parameter of the Green-Schwarz counterterm.  $\gamma_j$  are the anomalous dimensions of the matter fields, the  $\gamma_j^i$  and  $\gamma_j^{km}$  are their gauge and Yukawa parts, respectively.  $C_i$ ,  $C_i^j$  are the quadratic Casimir operators for the gauge group  $G_i$ , respectively, in the adjoint representation and in the matter representation. The indices i, j, k denote  $H_1$ ,  $H_2$ ,  $\tilde{E}$ ,  $\tilde{L}$ ,  $\tilde{D}$ ,  $\tilde{U}$  and  $\tilde{Q}$ . The A-parameters are finally given by:

$$A_{e,n}(GUT) = Y_{e,nn}(GUT)A_{\tilde{E}} \tilde{I}_{LH}, \tag{35}$$

$$A_{d,n}(GUT) = Y_{d,nn}(GUT)A_{\tilde{D}_n\tilde{Q}_nH_1}$$
(36)

$$A_{u,n}(GUT) = Y_{u,nn}(GUT)A_{\tilde{U}_n\tilde{Q}_nH_2}$$
(37)

where n denotes the generation.

In case of the OII compactification scheme the gaugino masses are given by Eqs. (3.11) of [33]:

$$M_{i} = -g_{i}^{2} m_{3/2} \left\{ \frac{\sqrt{3} \sin \theta}{2k_{s\bar{s}}^{1/2}} + \frac{1}{16\pi^{2}} \left[ 2t \cos \theta G_{2} \left( \delta_{GS} + b_{i} \right) + b_{i} + \frac{\sqrt{3}g_{s}^{2} \sin \theta}{2k_{s\bar{s}}^{1/2}} \left( C_{i} - \sum_{j} C_{i}^{j} \right) \right] \right\}.$$
(38)

For the sfermion parameters we have implemented two sets of boundary conditions: set (A) is specified by formulas Eqs. (3.15) and (3.19) of [33]:

$$M_i^2 = m_{3/2}^2 \left\{ \sin^2 \theta + \gamma_i + \frac{\sqrt{3} \sin \theta}{k_{c\bar{z}}^{1/2}} \left[ \sum_a \gamma_i^a g_a^2 + \frac{1}{2} \sum_{jk} \gamma_i^{jk} (k_s + k_{\bar{s}}) \right] \right\}, \tag{39}$$

$$A_{ijk} = m_{3/2} \left\{ \gamma_i + \gamma_j + \gamma_k - \frac{\sqrt{3}k_s \sin \theta}{k_{s\bar{s}}^{1/2}} \right\}$$

$$\tag{40}$$

Set (B) is specified by formulas Eqs. (3.16) and (3.20) of [33]:

$$M_{i}^{2} = m_{3/2}^{2} \left\{ \frac{\sqrt{3} \sin \theta}{k_{s\overline{s}}^{1/2}} \left[ 1 + 2t \cos \theta G_{2} \right] \left[ \sum_{a} g_{a}^{2} \gamma_{i}^{a} + \frac{1}{2} \sum_{jk} \gamma_{i}^{jk} \left( k_{s} + k_{\overline{s}} \right) \right] \right.$$

$$\left. + \sin^{2} \theta \left[ 1 + \gamma_{i} + \ln \left[ 2t | \eta(t)|^{4} \right] \left( \sum_{a} \gamma_{i}^{a} + 2 \sum_{jk} \gamma_{i}^{jk} \right) - \sum_{a} \gamma_{i}^{a} \ln(g_{a}^{2}) \right] \right.$$

$$\left. - \frac{9 \sin^{2} \theta}{k_{s\overline{s}}} \left[ \sum_{a} \gamma_{i}^{a} \left( \frac{g_{a}^{4}}{4} \right) \left( \ln(g_{a}^{2}) + \frac{5}{3} \right) \right.$$

$$\left. + \ln \left[ (t + \overline{t}) | \eta(t)|^{4} \right] \left( \sum_{a} \gamma_{i}^{a} \left( \frac{g_{a}^{4}}{4} \right) + \frac{1}{3} \sum_{jk} \gamma_{i}^{jk} k_{s} k_{\overline{s}} \right) \right] \right\}, \tag{41}$$

$$A_{ijk} = m_{3/2} \left\{ \left( \gamma_{i} + \gamma_{j} + \gamma_{k} \right) \left[ 1 + 2t \cos \theta G_{2} \right] \right.$$

$$\left. + \frac{\sqrt{3} \sin \theta}{k_{s\overline{s}}^{1/2}} \left[ k_{s} + \sum_{a} \frac{g_{a}^{2}}{2} \left( \gamma_{i}^{a} + \gamma_{j}^{a} + \gamma_{k}^{a} \right) \left( 1 - \ln(g_{a}^{2}) \right) \right.$$

$$\left. - \ln \left[ \left( t + \overline{t} \right) | \eta(t)|^{4} \right] \left( \sum_{a} g_{a}^{2} \left( \gamma_{i}^{a} + \gamma_{j}^{a} + \gamma_{k}^{a} \right) - \sum_{lm} k_{s} \left( \gamma_{i}^{lm} + \gamma_{j}^{lm} + \gamma_{k}^{lm} \right) \right) \right] \right\}. \tag{42}$$

In all three cases we have assumed that terms proportional to the  $\log(\tilde{\mu}_i)$  can be neglected  $(\tilde{\mu}_i)$  denote the Pauli Villar masses).

## 2.3.6 General High Scale Model

It is clear from the examples above that up to now there is no unique mechanism for supersymmetry breaking. Therefore, we have implemented the possibility to specify rather freely a high scale model. This model is specified by: a set of three in principal non–universal gaugino mass parameters  $M_{1/2}[U(1)]$ ,  $M_{1/2}[SU(2)]$ ,  $M_{1/2}[SU(3)]$ ; a scalar mass for each type of sfermion, resulting in fifteen parameters:  $M_{\tilde{E},ii}^0$ ,  $M_{\tilde{L},ii}^0$ ,  $M_{\tilde{D},ii}^0$ ,  $M_{\tilde{Q},ii}^0$ , two Higgs mass parameters  $M_{H_1}^0$  and  $M_{H_2}^0$ ; nine different A parameters  $A_{0,e,ii}$ ,  $A_{0,d,ii}$  and  $A_{0,u,ii}$ . Here ii denotes that only the diagonal entries can be set, because in the current version the effects of generation mixing is not taken into account. A model of this kind has been used in [34] for the study of low energy observables and the supersymmetric spectrum. It also can be used, for example, to set the boundary conditions for the gaugino mediated supersymmetry breaking [35]. This general model will be denoted by SUGRA.

## 2.3.7 General MSSM at low energies

Starting with version 2.2.0 there exists also the possibility to give the parameters at the low scale  $M_{EWSB}$  together with scale. In this case the parameters are taken as running parameters at the scale  $M_{EWSB}$  and the masses and mixing angles are calculated using these parameters which in turn serve as input for the calculation of decay widths and cross sections. The input parameter are: the electroweak gaugino mass parameters  $M_1$  and  $M_2$ , the gluino mass  $m_{\tilde{g}}$ ; the parameters describing the Higgs sector  $\mu$ , tan  $\beta$ , the mass of the pseudoscalar Higgs mass  $m_A$ ; the sfermion mass parameters  $M_{\tilde{E},ii}$ ,  $M_{\tilde{L},ii}$ ,  $M_{\tilde{D},ii}$ ,  $M_{\tilde{U},ii}$ ,  $M_{\tilde{Q},ii}$ ,  $A_{u,ii}$ ,  $A_{d,ii}$ ,

and  $A_{l,ii}$ . There are two general model implemented denoted by MSSM and pMSSM. The difference is in the interpretation of the parameters with respect to high order corrections as explained in Appendix B.3.8.

#### 3 Decays of supersymmetric particles and Higgs bosons

The programs calculates the most important two- and three-body decays of supersymmetric particles at tree level. In case of three-body decays the formulas are implemented such, that the effects of decay widths in the propagators are taken into account [36]. Therefore, it is possible to perform the calculation even in case that some of the intermediate particles are on-shell. This is useful in the case that the two-body decays have small phase space, because then the calculation of the three-body decays gives a more accurate result, e.g.  $\Gamma(\tilde{\chi}_1^+ \to 1)$  $\tilde{\chi}_1^0 W^+$ )×BR( $W^+ \to \nu l^+$ ) can be quite different from  $\Gamma(\tilde{\chi}_1^+ \to \tilde{\chi}_1^0 \nu l^+)$  if the decay  $\tilde{\chi}_1^+ \to \tilde{\chi}_1^0 W^+$ has only small phase space.

The following sfermion decays are calculated:

$$\tilde{f}_i \rightarrow f \, \tilde{\chi}_k^0, \, f' \, \tilde{\chi}_l^{\pm}$$
 (43)

$$\tilde{f}_i \rightarrow f \, \tilde{\chi}_k^0, \, f' \, \tilde{\chi}_l^{\pm} 
\tilde{f}_i \rightarrow \tilde{f}_j \, Z^0, \, \tilde{f}_j' W^{\pm}$$
(43)

$$\tilde{f}_i \rightarrow \tilde{f}_j (h^0, H^0, A^0), \ \tilde{f}'_j W^{\pm}$$
 (45)

In case of the lighter stop, it is possible that all two-body decays modes are kinematically forbidden at tree-level. In this case the following decay modes are important [37, 38, 39]:

$$\tilde{t}_1 \rightarrow c \, \tilde{\chi}_{1,2}^0 \tag{46}$$

$$\tilde{t}_1 \rightarrow W^+ b \, \tilde{\chi}_1^0, \ H^+ b \, \tilde{\chi}_1^0$$
 (47)

$$\tilde{t}_1 \rightarrow b \nu \tilde{l}_i^+, b l^+ \tilde{\nu}$$
 (48)

where  $l = e, \mu, \tau$ . The corresponding widths are calculated within SPheno using the formulas given in [38]. In case of GMSB models scenarios exist where the charged sleptons are next to lightest supersymmetric particles (NLSP) and the gravitino  $\hat{G}$  is the LSP. In this case the sleptons decay according to:

$$\tilde{l}_i \rightarrow l \, \tilde{G}$$
 (49)

Here we use the formulas given in [29].

It is well known that the partial widths of sfermions can receive considerable radiative corrections [40]. However, the branching ratios are not that strongly affected [41]. Therefore, for the moment being tree-level formulas are implemented. Some important numerical effects of higher order corrections are nevertheless implemented by using 1-loop corrected masses and running couplings in the formulas. The complete implementation of higher-order corrections is left for future versions of the program.

In case of charginos and neutralinos the following decay modes are calculated:

$$\tilde{\chi}_i^0 \rightarrow Z^0 \tilde{\chi}_j^0, W^{\pm} \tilde{\chi}_k^{\mp}$$
 (50)

$$\tilde{\chi}_{i}^{0} \rightarrow (h^{0}, H^{0}, A^{0}) \, \tilde{\chi}_{j}^{0}, \, H^{\pm} \, \tilde{\chi}_{k}^{\mp}$$
 (51)

$$\tilde{\chi}_i^0 \rightarrow f \bar{\tilde{f}}_j, \ \bar{f} \tilde{f}_j$$
 (52)

$$\tilde{\chi}_k^+ \rightarrow Z^0 \tilde{\chi}_s^+, W^+ \tilde{\chi}_j^0$$
 (53)

$$\tilde{\chi}_{k}^{+} \rightarrow (h^{0}, H^{0}, A^{0}) \, \tilde{\chi}_{s}^{+}, \, H^{+} \, \tilde{\chi}_{j}^{0}$$
 (54)

$$\tilde{\chi}_k^+ \to f \tilde{f}_i'$$
 (55)

In case that all two body decay modes are kinematically forbidden the following three-body decays are calculated:

$$\tilde{\chi}_i^0 \rightarrow f \bar{f} \tilde{\chi}_j^0, f f' \tilde{\chi}_k^{\mp}$$
 (56)

$$\tilde{\chi}_{i}^{0} \rightarrow q \bar{q} \tilde{g} \qquad (57)$$

$$\tilde{\chi}_{k}^{+} \rightarrow f \bar{f} \tilde{\chi}_{s}^{+}, f f' \tilde{\chi}_{j}^{0} \qquad (58)$$

$$\tilde{\chi}_{k}^{+} \rightarrow q q' \tilde{g} \qquad (59)$$

$$\tilde{\chi}_k^+ \rightarrow f \bar{f} \tilde{\chi}_s^+, f f' \tilde{\chi}_i^0$$
 (58)

$$\tilde{\chi}_k^+ \rightarrow q \, q' \, \tilde{g}$$
 (59)

In the calculation we have included all contributions from gauge bosons, sfermions and Higgs bosons [36, 42]. The Higgs contributions can be important in certain regions of parameter space [43]. In addition the loop induced decay s

$$\tilde{\chi}_i^0 \rightarrow \tilde{\chi}_i^0 \gamma$$
 (60)

are calculated [44] taking into account the left-right mixing of sfermions. Similarly to case of the sleptons there exist parameter regions in GMSB models where the lightest neutralino is the NLSP and it decays according to

$$\tilde{\chi}_1^0 \rightarrow \gamma \tilde{G}$$
 (61)

$$\tilde{\chi}_1^0 \rightarrow Z^0 \tilde{G}$$
 (62)

$$\tilde{\chi}_{1}^{0} \rightarrow \gamma \tilde{G} \qquad (61)$$

$$\tilde{\chi}_{1}^{0} \rightarrow Z^{0} \tilde{G} \qquad (62)$$

$$\tilde{\chi}_{1}^{0} \rightarrow h^{0} \tilde{G} \qquad (63)$$

Here we use the formulas given in [29].

In case of gluinos the following two-body decays are calculated:

$$\tilde{g} \to q \, \tilde{q}_i$$
 (64)

with q = u, d, c, s, t, b. Again, in case that these decays are kinematically suppressed, the three-body decay modes are calculated:

$$\tilde{g} \rightarrow \tilde{\chi}_i^0 q \bar{q}$$
 (65)

$$\tilde{g} \rightarrow \tilde{\chi}_j^{\pm} q' \bar{q}$$
 (66)

$$\tilde{g} \rightarrow \bar{b}W^{-}\tilde{t}_{1}, bW^{+}\tilde{t}_{1}^{*}$$

$$(67)$$

Here we have implemented the formulas given in [45]. In addition the decays

$$\tilde{g} \rightarrow \tilde{\chi}_i^0 g$$
 (68)

are calculated [44, 46] taking into account the left-right mixing of sfermions.

In case of Higgs bosons the following decays are calculated:

$$\phi \rightarrow f\bar{f} \tag{69}$$

$$\phi \rightarrow \tilde{f}_i \tilde{f}_j \tag{70}$$

$$\phi \rightarrow \tilde{f}_i \bar{\tilde{f}}_j 
\phi \rightarrow \tilde{\chi}_k^0 \tilde{\chi}_l^0$$
(70)

$$\phi \to \tilde{\chi}_r^+ \tilde{\chi}_s^- 
h^0 \to Z^0 Z^{0*}, W^+ W^{-*}, W^{+*} W^-$$
(72)

$$h^0 \rightarrow Z^0 Z^{0*}, W^+ W^{-*}, W^{+*} W^-$$
 (73)

$$h^0 \rightarrow gg \tag{74}$$

$$H^0 \rightarrow gg \tag{75}$$

$$H^0 \to Z^0 Z^0, W^+ W^-$$
 (76)

$$H^0 \rightarrow h^0 h^0 \tag{77}$$

$$A^0 \rightarrow h^0 Z^0 \tag{78}$$

$$H^+ \rightarrow f \bar{f}'$$
 (79)

$$H^+ \rightarrow \tilde{f}_i \, \bar{\tilde{f}}'_j$$
 (80)

$$H^{+} \rightarrow \tilde{\chi}_{k}^{0} \tilde{\chi}_{s}^{-}$$

$$H^{+} \rightarrow h^{0} W^{+}$$

$$(81)$$

$$H^+ \rightarrow h^0 W^+ \tag{82}$$

with  $\phi = h^0, H^0, A^0$  and  $f = \nu_i, e, \mu, \tau, u, d, c, s, t, b$ . It is well known, that the widths as well as the branching ratios of the Higgs bosons can receive large one-loop corrections [47, 48, 49]. In the present version only the gluonic QCD corrections for the decays into quarks [47] have been implemented. The decays into the gg final state have been implemented using the lowest order formula as given in [50]. Therefore, the numbers provided by SPheno have to be taken with care and for refined analysis other programs, such as HDECAY [51] should be used.

## Production of supersymmetric particles and Higgs 4 bosons

The program calculates the following cross sections:

$$e^+e^- \rightarrow \tilde{f}_i \tilde{f}_i \qquad (f=l,\nu,q)$$
 (83)

$$e^+ e^- \rightarrow \tilde{\chi}_r^+ \tilde{\chi}_s^-$$
 (85)

$$e^+e^- \rightarrow h^0 Z, H^0 Z$$
 (86)

$$e^{+}e^{-} \rightarrow h^{0}A^{0}, H^{0}A^{0}$$
 (87)

$$e^+e^- \rightarrow H^+H^-$$
 (88)

We haven taken the formulas of [52] for sfermion production, [53, 36] for production of charginos and neutralinos and [49] for Higgs boson production. Initial state radiation has been included using the formula given in [54]. In case of squarks in addition QCD corrections due to gluon exchange are included [54, 55]. Care has to be taken in case one calculates the cross sections near threshold because then higher order corrections are important to get reliable results [56] and, thus, the numbers obtained in the program have to be taken with care near the threshold. All cross sections are implemented such, that one can specify the degree of longitudinal polarization  $P_{e^-}$  of the incoming electron beam as well as the degree of longitudinal polarization  $P_{e^+}$  of the incoming positron beam. Here  $P_{e^-}$  is within the range [-1, 1], where  $\{-1, 0, 1\}$  denote 100% left-handed electrons, completely unpolarized electrons and 100% right-handed electrons, respectively. The same notation is used in case of positrons.

For example,  $P_{e^-} = -0.8$  ( $P_{e^+} = -0.8$ ) means that 80% of the electrons (positrons) are left-polarized whereas the remaining 20% are unpolarized.

## 5 Low Energy Constraints

The supersymmetric parameters are constrained by direct searches at colliders and by loop-effects which supersymmetric particles induce observables of low energy experiments. Provided one neglects mixing between different sfermion generations the following quantities constrain several parameters of the MSSM: the rare decay  $b \to s \gamma$ , the anomalous magnetic moment of the muon  $a_{\mu}$  and the supersymmetric contributions to the  $\rho$  parameter. These constraints are implemented in the program using the formulas given in [57, 58] for  $b \to s \gamma$  supplemented by the QCD corrections as given in [59], [60] for  $a_{\mu}$  and [61] for the sfermion contributions to the  $\rho$  parameter. In call cases we use the running couplings at  $m_Z$  for the calculation of the observables. The use of running couplings together with the correct implementation of supersymmetric threshold corrections for the couplings results in taking into account the most important higher oder corrections as has been pointed out e.g. in [62, 63] for the case of  $b \to s \gamma$ . The implementation of the supersymmetric threshold corrections to the couplings will be discussed in the next section.

## 6 Details of the Calculation

In this section we describe the algorithm used. It is schematically displayed in Fig. 1. The following standard model parameters are used as input: fermion masses, the Z-boson pole mass, the fine structure constant  $\alpha$ , the Fermi constant  $G_F$  and the strong coupling constant  $\alpha_s(m_Z)$ . It is assumed that  $\alpha_s(m_Z)$  is given in the  $\overline{MS}$  scheme. We describe first the implementation of the high scale models and comment then on the case of the implemented MSSM model.

## 6.1 First rough calculation of SUSY and Higgs boson masses

In a first step, we calculate gauge and Yukawa couplings at  $m_Z$  scale using tree-level formulas. These are used as input for the one-loop RGEs to get the gauge and Yukawa couplings at the high scale where also the boundary conditions for the high scale model under study are imposed. Afterward one-loop RGEs are used to get a first set of parameters at the electroweak scale. These parameters are used to get a first set of supersymmetric particle masses and Higgs masses using tree-level formulas except for the neutral CP-even Higgs bosons where one-loop effects due to (s)quarks of the third generation are taken into account. These masses are the starting point for the iterative loop which calculates the spectrum within the required precision as described below.

## 6.2 Main loop for the calculation of SUSY and Higgs boson masses

In the next step the gauge couplings and  $\sin^2 \theta_W$  are calculated at  $m_Z^2$  in the  $\overline{DR}$  scheme using the formulas given Appendix C of [25]. The only difference is that we use an updated number for  $\Delta \alpha_{lep} + \Delta \alpha_{had} = 0.06907$  [64]. The mass of the W-boson is calculated using the

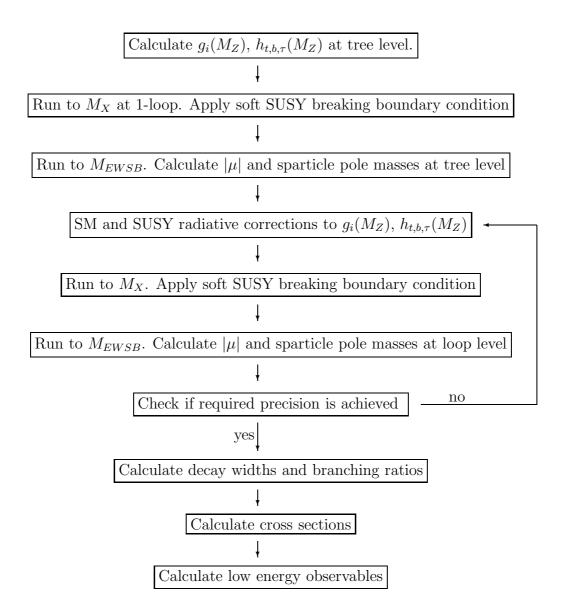


Figure 1: Algorithm used to calculate the SUSY spectrum, decay widths, production cross sections and low energy observables. Each step (represented by a box) is explained in the text. The initial step is the uppermost one.  $M_{EWSB}$  is the scale at which the EWSB conditions are imposed, and  $M_X$  is the scale at which the high energy SUSY breaking boundary conditions are imposed.

formula [65]

$$m_W^2 = m_Z^2 \hat{\rho} \left( \frac{1}{2} + \sqrt{\frac{1}{4} - \frac{\alpha^{\overline{DR}}(m_Z)\pi}{\sqrt{2}G_F m_Z^2 \hat{\rho}(1 - \Delta \hat{r})}} \right)$$
 (89)

$$\hat{\rho} = \frac{1}{1 - \Delta \hat{\rho}} \tag{90}$$

$$\Delta \hat{\rho} = Re \left( \frac{\Pi_{ZZ}^T(m_Z^2)}{\hat{\rho}m_Z^2} - \frac{\Pi_{WW}^T(m_W^2)}{m_W^2} \right)$$
(91)

$$\Delta \hat{r} = \hat{\rho} \frac{\Pi_{WW}^{T}(0)}{m_{W}^{2}} - \frac{\Pi_{ZZ}^{T}(m_{Z}^{2})}{\hat{m}_{Z}^{2}} + \delta_{VB}$$
(92)

where  $\Pi_{VV}^T$  is the transverse part of the vector boson self-energy and  $\delta_{VB}$  contains non-universal corrections to the  $\mu$  decay. In the computation we have included the leading SM two-loop corrections the formulas of [66] and adapting the Higgs contribution as in [25]. The formulas of the SUSY contribution to  $\delta_{VB}$  are given in [67, 68]. All masses appearing in the loops are running except for the top-quark, because the 2-loop part is given for an on-shell definition of the top mass. Note here that also the gauge boson masses in the loops are running masses and thus an iteration has to be performed in practice.

For the calculation of the Yukawa couplings we use the complete formulas for the fermion masses and the vacuum expectation values given in Appendix D of [25]. In addition we have implemented the following improvements. The five light quarks and  $\alpha_s$  are evolved to  $m_Z$  using 4-loop RGEs as given in [69] including threshold corrections at the low scale [70]. This evolution is done in the  $\overline{MS}$  scheme. At  $m_Z$  the shift to the  $\overline{DR}$  scheme is performed using the formulas given in [71, 72], e.g.:

$$m_{b',\overline{DR}}(m_Z) = m_{b,\overline{MS}}(m_Z) \left(1 - \frac{\alpha_s}{3\pi} - \frac{23\alpha_s^2}{72\pi^2} + \frac{3g_2^2}{128\pi^2} - \frac{13g'^2}{1152\pi^2}\right)$$
 (93)

where  $\alpha_s$  is given in the  $\overline{DR}$  scheme which is the reason for the different factor in front of  $\alpha_s^2$  compared to [72]. We use the complete formulas given in Appendix D of [25] to calculate the SUSY contribution denoted by  $\Delta m_{b,SUSY}(m_Z)$ . For the calculation we use running gauge and Yukawa couplings at  $m_Z$ . The SUSY contributions  $\Delta m_{b,SUSY}(m_Z)$  are then resumed using [73]

$$m_{b,\overline{DR}}(m_Z) = \frac{m_{b',\overline{DR}}(m_Z)}{1 - \frac{\Delta m_{b,SUSY}(m_Z)}{m_{b,\overline{DR}}(m_Z)}}$$
(94)

In the case of the top-quark a modified procedure has to be used due to its large mass. In this case we start from the pole mass  $m_t$  and use the formulas given in [25] to obtain  $m_t^{\overline{DR}}(m_Z)$ . The difference compared to [25] is in the 2-loop parts where we have taken the exact 2-loop formula given in [71]:

$$\Sigma_t^{2l} = -\left(\frac{\alpha_s(m_Z)}{\pi}\right)^2 \left(\frac{8\pi^2}{9} + \frac{2011}{18} + \frac{16\pi^2 \ln 2}{9} - \frac{8\zeta(3)}{3} + \frac{246}{3}L + 22L^2\right)$$
(95)

$$L = \log\left(\frac{m_Z^2}{m_t^2}\right) \tag{96}$$

In the case of leptons we first calculate the running masses  $m_l(m_l)$  at 1-loop level in the  $\overline{MS}$  scheme. Afterward we use two-loop RGEs as given in [74] to evolve them to  $m_Z$ . Here we perform the shift to the  $\overline{DR}$  scheme using the formula:

$$m_l^{\overline{DR},SM}(m_Z) = m_l^{\overline{MS},SM}(m_Z) \left(1 - \frac{3(g'^2 - g_2^2)}{128\pi^2}\right)$$
 (97)

The analog of Eq. (94) is used to obtain the running mass including the SUSY threshold corrections.

The obtained gauge and Yukawa couplings are evolved to the high scale using two-loop RGEs [24]. The high scale can either be fixed or can be calculated from the requirement  $g_1 = g_2$  at the high scale. The various cases are discussed in Sect. 7 and Appendix A. At the high scale the boundary conditions for the soft SUSY breaking parameters are set. The implemented models are summarized in Sect. 2.3; see also Appendix B.3. The complete set of two-loop RGEs [24] is used to evolve the parameters down to the electroweak scale  $M_{EWSB} = \sqrt{m_{\tilde{t}_1} m_{\tilde{t}_2}}$  (there exists also the possibility to fix  $M_{EWSB}$  to a constant value as described in Appendix A).

The parameters are used as input to calculate the sparticle pole masses at one-loop order and in case of the neutral Higgs at two-loop order. Here we use the complete formulae given in the appendices of [25] for the one-loop contributions and for the 2-loop corrections  $O(\alpha_s(\alpha_t + \alpha_b) + (\alpha_t + \alpha_b)^2 + \alpha_\tau \alpha_b + \alpha_t^2 au)$  for the neutral Higgs boson the formulas given in [26, 27, 75]. For the  $O(\alpha_s \alpha_b)$  contributions we use the complete expressions which can be obtained from the  $O(\alpha_s \alpha_t)$  contributions by appropriate replacements. In case of sfermions we have included in all cases left-right mixing. All gauge and Yukawa couplings are understood as  $\overline{DR}$  quantities at  $M_{EWSB}$ . Also  $\tan \beta$  and the vacuum expectation values are evaluated at  $M_{EWSB}$  to get a consistent set of input parameters. Note, that we express in all couplings the fermion masses and gauge boson masses by their corresponding expressions due to gauge couplings, Yukawa couplings and vacuum expectations values in the formulas of [25]. In all cases running masses are used as input for the loop integrals. In addition we have implemented the  $O(\alpha_s(\alpha_t + \alpha_b) + (\alpha_t + \alpha_b)^2 + \alpha_\tau \alpha_b + \alpha_t^2 au)$  corrections for the calculation of  $|\mu|$  [27, 75]. The numerical evolution of the one-loop integrals is based on the FF package [76] and the LoopTools package [77].

The masses obtained are used as input for the next iteration which starts again by calculating the SUSY contributions to gauge and Yukawa couplings at  $m_Z$ . In the case that during this iterative process an unphysical situation occurs, e.g. a pole mass squared being negative, the program terminates and it provides information on the exact reason for termination. The iteration is continued until all relative differences between the sparticle masses are smaller then the user imposed quantity  $\delta$ :

$$\delta > \frac{|m_i - m_{i-1}|}{m_i} \tag{98}$$

for all sparticle masses; i denotes the i-th iteration. In most cases this achieved after three to four iterations. In the case that more than the maximal allowed number of iterations (user specified) are necessary, the program leaves the iteration giving a warning message.

In the case of the MSSM model the running between  $M_{EWSB}$  and  $M_X$  is obviously omitted. In this model one has, however, to preform an iteration to get the Yukawa and gauge couplings correct because  $\tan \beta$  is defined at  $M_{EWSB}$  instead of  $m_Z$ .

## 6.3 Calculation of the other observables

The masses and mixing angles are then used to calculate the branching ratios and decay widths. Here two- and three body decays of supersymmetric particles are calculate. Note that we use the couplings as input which are renormalized at the scale  $M_{EWSB}$ . The user has the possibility to force the program to calculate three body decays even if one or more of the intermediate particles are on-shell. This possibility is useful in the case where the 2-body decay has only small phase because then the calculation of the three body decay width(s) give a more reliable result.

Afterward the production cross sections at an  $e^+e^-$  collider of all supersymmetric particles as well as all Higgs bosons are calculated. Here the user has the possibility to specify the center of mass energy as well as the degree of longitudinal polarization of the incoming beams. Moreover, the user can specify if initial state radiation shall be included in the calculation or not.

Finally, the low–energy constraints described in Section 5 are calculated:  $BR(b \to s\gamma)$ , SUSY contributions to  $a_{\mu}$  and the sfermion contributions to  $\Delta \rho$ . For the the calculation of these quantities we evolve the gauge and Yukawa couplings from the scale  $M_{EWSB}$  down to  $m_Z$ . The couplings at  $m_Z$  are then used as input for the calculation of the low energy observables. For example in calculation of  $BR(b \to s\gamma)$  the most important contributions to the  $C_7$  coefficients are implemented as

$$C_{7}(W^{+}) = -\frac{K_{ts}K_{tb}x_{tW}}{4 m_{W}^{2}} \left(\frac{2}{3}F_{1}(x_{tW}) + F_{2}(x_{tW})\right)$$

$$C_{7}(H^{+}) = -\frac{K_{ts}K_{tb}}{4 m_{H^{+}}^{2}} \left[\frac{Y_{t}^{2}\cos^{2}\beta}{4} \left(\frac{2}{3}F_{1}(x_{tH^{+}}) + F_{2}(x_{tH^{+}})\right) - \frac{Y_{b}Y_{t}\cos\beta\sin\beta m_{t}}{m_{b}} \left(\frac{2}{3}F_{3}(x_{tH^{+}}) + F_{4}(x_{tH^{+}})\right)\right]$$

$$C_{7}(\tilde{\chi}^{+}) = \sum_{i,j=1}^{2} \frac{K_{ts}K_{tb}}{4 m_{\tilde{t}_{i}}^{2}} \left[C_{R,ij}^{2} \left(\frac{2}{3}F_{2}(x_{\tilde{\chi}_{j}^{+}\tilde{t}_{i}}) + F_{1}(x_{\tilde{\chi}_{j}^{+}\tilde{t}_{i}})\right) \right]$$

$$(99)$$

$$-C_{L,ij}C_{R,ij}\left(\frac{2}{3}F_4(x_{\tilde{\chi}_i^+\tilde{t}_i}) + F_3(x_{\tilde{\chi}_i^+\tilde{t}_i})\right)\right]$$
(101)

$$C_{L,ij} = Y_b R_{\tilde{t},i1} U_{j2} \tag{102}$$

$$C_{R,ij} = -gR_{\tilde{t},i1}V_{j1} + Y_tR_{\tilde{t},i2}V_{j2} \tag{103}$$

Here  $Y_i$  are the Yukawa couplings, U and V are the chargino matrices, K is the CKM matrix,  $R_{\tilde{t}}$  is the stop mixing matrix and  $x_{ab} = m_a^2/m_b^2$ . The loop functions  $F_i$  are given in [57]. A similar replacement is done in the contributions to the  $C_8$  coefficient. Moreover, in the program also the contributions from the first two generation of (s)fermions to  $C_{7,8}$  are included for completeness. We then use [59] to obtain

$$BR(b \to s\gamma) = 1.258 + 0.382r7^2 + 0.015r_8^2 + 1.395r_7 + 0.161r_8 + 0.083r_7r_8$$
 (104)

where  $r_7 = C_7/C_7(W^+)$  and  $r_8 = C_8/C_8(W^+)$ . In this way important higher order corrections are taken into account, in particular the large  $\tan \beta$  effects in case of  $b \to s\gamma$  [63].

Table 1: Variables for parameters and couplings. The parameters are defined in the module MSSM\_Data and they are explained in Section 2. dp means double precision.

```
parameter/coupling
                     type & Fortran name
e^{\varphi_{\mu}}
                     complex(dp) :: phase_mu
\tan \beta
                     real(dp) :: tanb
M_1, M_2, M_3
                     complex(dp) :: Mi(3)
M_E^2, M_L^2
                     complex(dp), dimension(3,3) :: M2_E, M2_L
M_D^2, M_Q^2, M_U^2
                     complex(dp), dimension(3,3) :: M2_D, M2_Q, M2_U
A_l, A_d, A_u
                     complex(dp), dimension(3,3) :: A_1, A_d, A_u
                     complex(dp) :: mu
\mu
B\mu
                     complex(dp) :: B
M_H^2
                     real(dp) :: M2_H(2)
g', g
                     real(dp) :: gp, g
Y_l, Y_d, Y_u
                     complex(dp), dimension(3,3) :: Y_1, Y_d, Y_u
                     real(dp) :: vevSM(2)
v_1, v_2
                     real(dp) :: gauge(3)
g', g, g_s
```

# 7 A sample example

In this section we discuss the executable statements of the main program given in the file SPheno.f90. In the first statements the required modules are loaded and the various variables are defined. Afterward the error system is initialized and the input data are read by calling:

## Call ReadingData(kont)

The source code can be found in the file SPheno.f90. This routine checks first if the file LesHouches.in exists where the input data are provided according to SLHA (for a detailed description see Appendix C). If this file is absent, the routine checks if the files Control.in and/or StandardModel.in are present. The first one can be used to set various flags (see Appendix B.1) whereas the second one is used to set the SM input (see Appendix B.4). Standard values as described in the Appendix are used if any of these files is absent. Afterward the model specific information is read from the file HighScale.in which is described in Appendix B.3 (for a short description of the implemented models see Sect. 2.3). In all cases the file Messages.out is opened at channel 10 where all warnings and/or debugging informations are stored.

Before calling the subroutine CalculateSpectrum, which performs the calculation of the spectrum, the user has the possibility to fix the high scale and/or the scale where the parameters and the loop corrected masses are calculated. For this purpose one or both of the following lines must be uncommented in the program:

```
! Call SetGUTScale(2.e16_dp) ! please put the GUT scale
! Call SetRGEScale(1.e3_dp**2) ! please put the scale M_EWSB squared
```

The default is that these scales are calculated by the program. The high scale is computed from the requirement  $g_1 = g_2$  (except in GMSB where the high scale is an input). The scale  $M_{EWSB}$  is given by  $M_{EWSB} = \sqrt{m_{\tilde{t}_1} m_{\tilde{t}_2}}$ .

The accurate calculation of the SUSY parameters and the spectrum is done by the following call:

Table 2: Variables for masses and mixing matrices as given by the routine CalculateSpectrum. They are defined in the module MSSM\_Data and their connection to the parameters at tree—level is explained explained in Section 2. dp means double precision.

```
masses / mixing matrix type & Fortran name
                                  real(dp) :: mglu
m_{\tilde{q}}
e^{\varphi_{\tilde{g}}}
                                  complex(dp) :: PhaseGlu
                                  real(dp) :: mC(2)
m_{\tilde{\chi}_i^+}
U, V
                                  complex(dp) :: U(2,2), V(2,2)
                                  real(dp) :: N(4)
m_{\tilde{\chi}^0_i}
N
                                  complex(dp) :: N(4,4)
                                  real(dp) :: mSO(2)
m_{h^0}, m_{H^0}
                                  real(dp) :: RSO(2,2)
R_{\alpha}
                                  real(dp) :: mP0(2)
m_{G^0}, m_{A^0}
                                  real(dp) :: RPO(2,2)
R_{\beta}
                                  real(dp) :: mSpm(2)
m_{G^+}, m_{H^+}
R'_{\beta}
                                  complex(dp) :: RSpm(2,2)
                                  real(dp) :: mSneut(3)
m_{\tilde{\nu}}
                                  complex(dp) :: Rsneut(3,3)
R_{\tilde{\nu}}
                                  real(dp) :: mSlepton(6)
m_{\tilde{i}}
                                  complex(dp) :: Rslepton(6,6)
R_{\tilde{l}}
                                  real(dp) :: mSup(6)
m_{\tilde{u}}
                                  complex(dp) :: Rsup(6,6)
R_{\tilde{u}}
                                  real(dp) :: mSdown(6)
m_{\tilde{d}}
R_{\tilde{d}}
                                  complex(dp) :: Rsdown(6,6)
```

```
delta = 1.e-4_dp
WriteOut = .False.
n_run = 20
If (kont.Eq.0) Call CalculateSpectrum(n_run, delta, WriteOut, kont, tanb & vevSM, mC, U, V, mN, N, mSO, mSO2, RSO, mPO, mPO2, RPO, mSpm, mSpm2 & RSpm, mSdown, mSdown2, RSdown, mSup, mSup2, RSup, mSlepton, mSlepton2 & RSlepton, mSneut, mSneut2, RSneut, mGlu, PhaseGlu, gauge, Y_l, Y_d & Y_u, Mi, A_l, A_d, A_u, M2_E, M2_L, M2_D, M2_Q, M2_U, M2_H, mu, B & A_l_save, A_u_save, A_d_save, m_GUT)
```

The meaning of the various variables and their type is given in Tables 1 and 2. Variable names ending with "2" indicate masses squared. The variables for the mixing matrices are already structured for a latter extension to include the effects of generation mixing and/or complex phases: the sfermion mixing matrices are  $6 \times 6$  (except for sneutrinos which is a  $3 \times 3$  matrix). In the present release most of the entries are zero except for the diagonal  $2 \times 2$  blocks which contain the left–right mixing for every species of sfermions. For example, the 11, 12, 21 and 22 entries in Rslepton specify the left–right mixing of selectrons, and similarly the 33, 34, 43 and 44 (55, 56, 65 and 66) entries specify the left–right mixing of smuons (staus). The squark mixing matrices have the same generation structure. This structure has been chosen to facilitate a later extension which includes flavour violating entries. Beside the variables given in Tables 1 and 2 the following variables appear:

- delta: specifies the required relative precision on the masses. If the maximal relative differences between the physical masses obtained between two runs is smaller than delta, the routine Sugra leaves the iteration loop.
- m\_GUT: the value of the scale where the high energy boundary conditions are imposed.
- kont: A variable which is 0 provided everything is o.k. Otherwise either a numerical problem has occurred and/or the parameters belong to an unphysical region, e.g. a minimum of the potential where charge and/or colour breaking minima occur. In such a case the information is written to the file Messages.out.
- WriteOut: If it is set .True. then intermediate debugging information is written to the screen and the file Messages.out.
- n\_run specifies the maximal number of iterations of the main loop. A warning will be given in the case that the required precision delta has not been reached within n\_run iterations.

Note that the parameters are running parameters at the scale  $M_{EWSB}$ . The complete spectrum is calculated at 1-loop level using the formulas given in [25]. The exceptions are the masses of the neutral Higgs bosons (scalar and pseudo-scalar) as well as  $\mu$  the two loop corrections due to  $\alpha_s$  and all third generation Yukawa couplings are included [26, 27].

In the next part the branching ratios, the partial decay widths and the total decay widths are calculated provided that L\_BR=.TRUE. and kont.eq.0:

```
If ((L_BR).And.(kont.eq.0)) Then
epsI = 1.e-5_dp
deltaM = 1.e-3_dp
CalcTBD = .False.
ratioWoM = 0._dp
Couplings_At_Mi = .True.
Call CalculateBR(gauge, mGlu, PhaseGlu, mC, U, V, mN, N, mSneut
                                                                             &
   & , mSlepton, RSlepton, mSup, RSup, mSdown, RSdown, mSO, RSO
                                                                             &
   & , mPO, RPO, mSpm, RSpm, Y_d, A_d, Y_1, A_1, Y_u, A_u, mu, vevSM
                                                                             &
   & , Fgmsb, m32 , epsI, deltaM, ratioWoM , CalcTBD, Couplings_At_Mi
                                                                             &
   & , kont, gP_Sl, gT_Sl, BR_Sl, gP_Sn, gT_Sn, BR_Sn, gP_Sd, gT_Sd, BR_Sd &
   & , gP_Su, gT_Su, BR_Su, gP_C, gT_C, BR_C, gP_N, gT_N, BR_N
                                                                             &
   & , gP_Glu, gT_Glu, BR_Glu, gP_PO, gT_PO, BR_PO, gP_SO, gT_SO, BR_SO
                                                                             &
   & , gP_Spm, gT_Spm, BR_Spm)
End If
```

Variables starting with gP\_, gT\_ and BR\_ indicate partial widths, total widths and branching ratios, respectively; they are Real(dp) vectors. The first index is the index of the decaying particle whereas the second one gives the mode. The correspondence between the second index and the modes is summarized for sfermions (variables ending S1, Sn, Sd and Su for sleptons, sneutrino, d-squarks and u-squarks, respectively) in Table 3, for charginos in Table 4, for neutralinos in Table 5, for gluinos in Table 6 and for the Higgs bosons in Tables 7 and 8.

Here the following variables are new:

Table 3: Correspondence between the indices for sfermion partial widths (branching ratios) and the modes.  $\tilde{t}_1$  is listed extra because it can have three-body decay modes.

mode	$\widetilde{l}$	$ ilde{ u}$	$ ilde{d}$	$ ilde{u}$	$ ilde{t}_1$
$\tilde{f}_i \to f \tilde{\chi}_k^0$	1-4	1-4	1-4	1-4	1-4
$\tilde{f}_i \to f' \tilde{\chi}_j^{\pm}$	5-6	5-6	5-6	5-6	5-6
$ ilde{f_i}  ightarrow f  ilde{g}$	-	-	7	7	7
$\tilde{f}_i \to W^{\pm} \tilde{f}'_i$	7	7-8	8-9	8-9	8-9
$\tilde{f}_i \to H^{\pm} \tilde{f}'_i$	8	9-10	10-11	10-11	10-11
$ ilde{f}_2  ightarrow Z^0  ilde{f}_1$	9	-	12	12	-
$\tilde{f}_2 \to A^0 \tilde{f}_1$	10	-	13	13	-
$\tilde{f}_2 \to h^0 \tilde{f}_1$	11	-	14	14	-
$ ilde{f}_2  o H^0  ilde{f}_1$	12	-	15	15	-
$\tilde{l}_1 \to l  \tilde{G}$	13	-	-	-	-
$ \begin{array}{c} \tilde{t}_1 \to c\tilde{\chi}^0_{1,2} \\ \tilde{t}_1 \to W^+ \bar{b}\tilde{\chi}^0_1 \end{array} $	-	-	-	-	55-56
$\tilde{t}_1 \to W^+ \bar{b} \tilde{\chi}_1^0$	-	-	-	-	57
$\tilde{t}_1 \to \bar{b}e^+\tilde{\nu}_e$	-	-	-	-	58
$\tilde{t}_1 \to \bar{b}\mu^+\tilde{\nu}_\mu$	-	-	-	-	59
$\tilde{t}_1 \to \bar{b} \tau^+ \tilde{\nu}_{\tau}$	-	-	-	-	60
$\tilde{t}_1 \to \bar{b}\nu_e \tilde{e}_{1,2}^+$	-	_	-	-	61-62
$\tilde{t}_1 \rightarrow \bar{b}\nu_{\mu}\tilde{\mu}_{1,2}^+$	_	-	-	-	63-64
$\tilde{t}_1 \to \bar{b}\nu_{\tau}\tilde{\tau}_{1,2}^+$	-	_	-	-	65-66

- Fgmsb and m32: the F parameter and the gravitino mass in the GMSB model. These parameters are calculated from the input and are set to huge numbers in all other models. They are needed for the calculation of the decay width(s) of the NLSP into a gravitino.
- epsI: gives the accuracy to which the phase space integrals in three body decays are calculated.
- deltaM: this variable affects the calculation of the phase space integrals in three body decays. In case that  $m_i/(m-\sum_i m_i) < \text{deltaM}$  than  $m_i$  is set to zero in the calculation of the phase space integrals. m denotes here mass of the decaying particle and  $m_i$  (i=1,2,3) are the masses of the decay products.
- ratioWoM: this variable is used to decide whether two body decays or three body decay modes are calculated in case of charginos, neutralino and gluino. The program tries first two-body decay modes. In the case that the ratio of the width  $\Gamma$  over the mass m of the decaying particle is small:  $\Gamma/m < \text{ratioWoM}$ , then three body decay modes are calculated.
- CalcTBD: if this variable is set .TRUE. then in all chargino-, neutralino- and gluino decays the three body modes will be calculated. This option has to be taken with care, because it can slow down the program considerably.

Table 4: Correspondence between the second indices for chargino partial widths (branching ratios) and the decay modes.

mode	index of gP_C (BR_C)			
$\tilde{\chi}_i^+ \to \tilde{l}_{m,k}^+  \nu_m$	1-6			
$\tilde{\chi}_i^+ \to \tilde{\nu}_m  l_m^+$	7-9			
$\tilde{\chi}_i^+ \to \bar{\tilde{d}}_{m,k} u_m$	10-15			
$\tilde{\chi}_i^+ \to \tilde{u}_{m,k}  \bar{d}_m$	16-21			
$\tilde{\chi}_i^+ \to \tilde{\chi}_i^0 W^+$	22-25			
$\tilde{\chi}_i^+ \to \tilde{\chi}_i^0 H^+$	26-29			
$\tilde{\chi}_2^+  ightarrow \tilde{\chi}_1^+ Z^0$	30			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ A^0$	31			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ h^0$	32			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ H^0$	33			
$\tilde{\chi}_i^+ \to \tilde{\chi}_j^0 u_m  \bar{d}_m$	64-75			
$\tilde{\chi}_i^+ \to \tilde{\chi}_j^0 l_m^+ \nu_m$	76-87			
$\tilde{\chi}_i^+ \to \tilde{g} u_m  \bar{d}_m$	88-90			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ u_m  \bar{u}_m$	91-93			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+  d_m  \bar{d}_m$	94-96			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ l_m l_m^+$	97-99			
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ \nu_m  \bar{\nu}_m$	100-102			

• Couplings\_At\_Mi: if this variable is set .TRUE. then all couplings will be evolved to the scale  $Q=M_i$ , where  $M_i$  is the mass of the decaying particle. If this variable is set .False. then the couplings at the scale  $Q=m_{EWSB}$  will be used.

The next statements call the routine for the calculation of the cross sections provided  $L_CS = .TRUE$ . and kont = 0:

End If

Here the following additional input is needed:

- Ecms: the center of mass energy of the collider
- Pm, Pp: degree of polarization of the incoming electron and positron, respectively
- ISR: logical variable, if .TRUE. then initial state radiation is taken into account using the formulas given in [54]

Table 5: Correspondence between the second indices for neutralino partial widths (branching ratios) and the decay modes. Note, that also charge conjugated states are given.

$\operatorname{mode}$	index of gP_N (BR_N)
$\tilde{\chi}_i^0 \to \tilde{l}_{m,k}^+  l_m$	1-12
$\tilde{\chi}_i^0 \to \tilde{\nu}_m  \nu_m$	13-18
$\tilde{\chi}_i^0 \to \bar{\tilde{u}}_{m,k} u_m$	19-30
$\tilde{\chi}_i^0 \to \tilde{d}_{m,k}  \bar{d}_m$	31-42
$\tilde{\chi}_i^0 \to \tilde{\chi}_i^{\pm} W^{\mp}$	43-46
$\begin{array}{c} \tilde{\chi}_{i}^{0} \rightarrow \tilde{\chi}_{j}^{\pm} W^{\mp} \\ \tilde{\chi}_{i}^{0} \rightarrow \tilde{\chi}_{j}^{\pm} H^{\mp} \end{array}$	47-50
$\begin{array}{c} \tilde{\chi}_{i}^{0} + \tilde{\chi}_{j}^{0} \Pi \\ \tilde{\chi}_{i}^{0} \to \tilde{\chi}_{j}^{0} Z^{0} \\ \tilde{\chi}_{i}^{0} \to \tilde{\chi}_{j}^{0} A^{0} \end{array}$	51-(24+i)
$\tilde{\chi}_i^0 \to \tilde{\chi}_i^0 A^0$	(25+i)-(23+2i)
$\tilde{v}^0 \rightarrow \tilde{v}^0 h^0$	(26+i)-(22+3 i)
$\tilde{\chi}_i^0  ightarrow \tilde{\chi}_j^0 H^0$	(27+i)-(21+4i)
$\frac{\tilde{\chi}_{i}^{0} \to \tilde{\chi}_{j}^{0} H^{0}}{\tilde{\chi}_{i}^{0} \to \gamma \tilde{G}}$	63
$\tilde{\chi}^0_i  o Z^0  \tilde{G}$	64
$ ilde{\chi}^0_i  ightarrow h^0   ilde{G}$	65
$\frac{\tilde{\chi}_i^0 \to h^0  \tilde{G}}{\tilde{\chi}_i^0 \to \gamma  \tilde{\chi}_j^0}$	65 + j
$\tilde{\chi}_i^0 \to \tilde{\chi}_i^{\pm} q_m  \bar{q}_m'$	69 - 80
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^{\pm} l_m^{\mp} \nu_m$	81 - 92
$\tilde{\chi}_i^0 \to \tilde{G} u_m  \bar{u}_m$	93 - 95
$\tilde{\chi}_i^0 \to G  d_m  \bar{d}_m$	96 - 98
$\tilde{\chi}_{i>1}^0 \to \tilde{\chi}_1^0 u_m  \bar{u}_m$	99 - 101
$\tilde{\chi}_{i>1}^0 \to \tilde{\chi}_1^0 d_m \bar{d}_m$	102 - 104
$\tilde{\chi}_{i>1}^0 \to \tilde{\chi}_{\frac{1}{2}}^0 l_m^+ l_m^-$	105 - 107
$\tilde{\chi}_{i>1}^0 \to \tilde{\chi}_1^0 \nu_m \bar{\nu}_m$	108 - 110
$\tilde{\chi}_{i>2}^0 \to \tilde{\chi}_2^0 u_m \bar{u}_m$	111 - 113
$\tilde{\chi}_{i>2}^0 \to \tilde{\chi}_2^0 d_m \bar{d}_m$	114 - 116
$\tilde{\chi}_{i>2}^{0} \to \tilde{\chi}_{2}^{0} l_{m}^{+} l_{m}^{-}$	117 - 119
$ \tilde{\chi}_{i>2}^{0} \to \tilde{\chi}_{2}^{0} \nu_{m} \bar{\nu}_{m}  \tilde{\chi}_{4}^{0} \to \tilde{\chi}_{3}^{0} u_{m} \bar{u}_{m} $	120 - 122
$\chi_4^0 \rightarrow \bar{\chi}_3^0 u_m \bar{u}_m$	123 - 125
$\tilde{\chi}_4^0 \rightarrow \tilde{\chi}_3^0 d_m \bar{d}_m$	126 - 128
$\tilde{\chi}^0_4 \rightarrow \tilde{\chi}^0_3 l_m^+ l_m^-$	129 - 131
$\tilde{\chi}_4^0 \to \tilde{\chi}_3^0  \nu_m  \bar{\nu}_m$	132 - 134

Table 6: Correspondence between the indices for gluino partial widths (branching ratios) and the decay modes. Note, that also charge conjugated states are given.

mode	index of gP_G (BR_G)
$\tilde{g} \to \bar{\tilde{d}}_{m,k} d_m$	1-12
$\tilde{g} \to \tilde{u}_{m,k}  \bar{u}_m$	13-24
$\tilde{g} \to \tilde{t}_1  \bar{c}$	25-26
$\tilde{g} \to \gamma  \tilde{\chi}_i^0$	26 + j
$\tilde{g} \to \tilde{\chi}_1^0  u_m  \bar{u}_m$	31-33
$\tilde{g} \to \tilde{\chi}_1^0  d_m  \bar{d}_m$	34-36
$\tilde{g} \to \tilde{\chi}_2^0  u_m  \bar{u}_m$	37-39
$\tilde{g} \to \tilde{\chi}_2^0  d_m  \bar{d}_m$	40-42
$\tilde{g} \to \tilde{\chi}_3^0  u_m  \bar{u}_m$	43-45
$\tilde{g} \rightarrow \tilde{\chi}_3^0  d_m  \bar{d}_m$	46-48
$\tilde{g} \to \tilde{\chi}_4^0 u_m  \bar{u}_m$	49-51
$\tilde{g} \to \tilde{\chi}_4^0  d_m  \bar{d}_m$	52-54
$\tilde{g} \to \tilde{\chi}_i^{\pm} q_m  \bar{q}_m'$	55-66
$\tilde{g} \to \tilde{t}_i W^- \bar{b}^{m}$	67-68

These variables can be set in the file CrossSections.in. The cross sections are stored in the variables starting with Sig which are summarized in Table 9. Please note, that in case of sfermions the structure of the variables is already put such that the case of generation mixing can easily be implemented. In the non-mixing case the cross sections are stored in the  $2 \times 2$  diagonal blocks and they are sorted according to the generations as in the case of the sfermion mixing matrices.

Finally the low energy constraints  $b \to s\gamma$ ,  $a_{\mu}$  and  $\Delta \rho$  are calculated provided that calculation of the spectrum had been performed successfully (kont.eq.0):

Here BRbtosgamma, a\_mu, and Delta\_Rho denote  $10^4 \times \text{BR}(b \to s\gamma)$ , the SUSY contributions to  $a_{\mu}$  and the sfermion contributions to  $\Delta \rho$ , respectively.

Afterward the output is written using the following sequence of commands:

```
Call SetWriteMinBR(1.e-3_dp) ! sets the minimum branching ratio, which is written Call SetWriteMinSig(1.e-3_dp) ! sets the minimum cross section, which is written
```

Table 7: Correspondence between the indices for the partial widths (branching ratios) of the neutral Higgs bosons and the decay modes. The variables are gP\_S0 (BR\_S0) and gP\_P0 (BR\_P0) for the partial decay widths (branching ratios) of the CP-even Higgs bosons ( $h^0$ ,  $H^0$ ) and CP-odd Higgs boson ( $A^0$ ). In case of gP\_S0 (BR\_S0) the first (second) index denotes decay modes of  $h^0$  ( $H^0$ ). Here  $\phi$  stands for  $h^0$ ,  $H^0$  and  $A^0$ . The index m runs from 1 to 3.

mode	$h^0$	$H^0$	$A^0$	mode	$h^0$	$H^0$	$A^0$
$\phi \to l_m^+ l_m^-$	1-3	1-3	1-3	$H^0  o Z^0 Z^0$	-	63	-
$\phi \to d_m  \bar{d}_m$	4-6	4-6	4-6	$H^0  o W^+ W^-$	-	64	-
$\frac{\phi \to u_m  \bar{u}_m}{H^0 \to \tilde{e}_1^+  \tilde{e}_1^-}$	7-9	7-9	7-9	$H^0 \to h^0 h^0$	-	70	-
	-	10	-	$A^0 \rightarrow h^0 Z^0$	-	-	63
$\phi \to \tilde{e}_1^{\mp}  \tilde{e}_2^{\pm}$	-	11-12	11-12	$h^0 \to W^+W^{-*}$	70	-	-
$H^0  o \tilde{e}_2^+  \tilde{e}_2^-$	-	13	-	$h^0 \to W^{+*}W^-$	71	-	-
$H^0  o \tilde{\mu}_1^+ \tilde{\mu}_1^-$	-	14	-	$h^0 \to Z^0 Z^{0^*}$	72	-	-
$\phi \to \tilde{\mu}_1^{\mp}  \tilde{\mu}_2^{\pm}$	-	15-16	15-16	$\phi \to gg$	80	80	-
$H^0  ightarrow  ilde{\mu}_2^+   ilde{\mu}_2^-$	-	17	-				
$H^0  o  ilde{ au}_1^+  ilde{ au}_1^-$	-	18	-				
$\phi \to \tilde{\tau}_1^{\mp}  \tilde{\tau}_2^{\pm}$	-	19-20	19-20				
$H^0  o  ilde{ au}_2^+  ilde{ au}_2^-$	-	21	-				
$H^0  o \tilde{\nu}_m  \overline{\tilde{\nu}}_m$	-	21 + m	-				
$H^0  ightarrow \tilde{d}_1^+  \tilde{d}_1^-$	-	25	-				
$\phi \to \tilde{d}_1^{\mp}  \tilde{d}_2^{\pm}$	-	26-27	23-24				
$H^0  ightarrow  ilde{d}_2^+  ilde{d}_2^-$	-	28	-				
$H^0  ightarrow \tilde{s}_1^+  \tilde{s}_1^-$	-	29	-				
$\phi \to \tilde{s}_1^{\mp}  \tilde{s}_2^{\pm}$	-	30-31	27-28				
$H^0  ightarrow \tilde{s}_2^+  \tilde{s}_2^-$	-	32	-				
$H^0  ightarrow  ilde{b}_1^+  ilde{b}_1^-$	-	33	-				
$\phi  o \tilde{b}_1^{\mp}  \tilde{b}_2^{\pm}$	-	34-35	31-32				
$H^0  ightarrow  ilde{b}_2^+   ilde{b}_2^-$	-	36	-				
$H^0  ightarrow \tilde{u}_1^+  \tilde{u}_1^-$	-	37	-				
$\phi \to \tilde{u}_1^{\mp}  \tilde{u}_2^{\pm}$	-	38-39	35-36				
$H^0  ightarrow  ilde{u}_2^+   ilde{u}_2^-$	-	40	-				
$H^0  o \tilde{c}_1^+  \tilde{c}_1^-$	-	41	-				
$\phi \to \tilde{c}_1^{\mp}  \tilde{c}_2^{\pm}$	-	42-43	39-40				
$H^0  ightarrow  ilde{c}_2^+   ilde{c}_2^-$	-	44	-				
$H^0  o \tilde{t}_1^+  \tilde{t}_1^-$	-	45	-				
$\phi \to \tilde{t}_1^{\mp} \tilde{t}_2^{\pm}$	-	46-47	43-44				
$\frac{H^0 \to \tilde{t}_2^+  \tilde{t}_2^-}{\phi \to \tilde{\chi}_r^0  \tilde{\chi}_s^0  (r \le s)}$	-	48	-				
$\phi \to \tilde{\chi}_r^0  \tilde{\chi}_s^0  \left( r \le s \right)$	49-58	49-58	46-55				
$\phi \to \tilde{\chi}_1^+ \tilde{\chi}_1^-$	59	59	56				
$\phi \to \tilde{\chi}_1^{\pm}  \tilde{\chi}_2^{\mp}$	60-61	60-61	57-58				
$\phi \to \tilde{\chi}_2^+  \tilde{\chi}_2^-$	62	62	59				

Table 8: Correspondence between the indices for the partial widths (branching ratios) of the charged Higgs and the decay modes. The variables are gP\_Spm (BR\_Spm) and in case of partial decay widths (branching ratios).

mode	index
$H^+ \to l_m^+ \nu_m$	1-3
$H^+ \to u_m  \bar{d}_m$	4-6
$H^+ \to \tilde{e}_i^+  \tilde{\nu}_e$	7-8
$H^+  o \tilde{\mu}_i^+  \tilde{\nu}_\mu$	9-10
$H^+ \to \tilde{ au}_i^+  \tilde{ au}_{ au}$	11-12
$H^+  o \tilde{u}_i  \overline{\tilde{d}}_j$	12 + 2*(j-1) + i
$H^+ \to \tilde{c}_i  \overline{\tilde{s}}_j$	16 + 2*(j-1) + i
$H^+  o \tilde{t}_i  \overline{\tilde{b}}_j$	20 + 2*(j-1) + i
$H^+ \to \tilde{\chi}_r^+  \tilde{\chi}_s^0$	24 + 4*(r-1) + s
$H^+ \to h^0 W^+$	34

Table 9: Correspondence between the production cross sections and the variables used in the program.

process	Fortran name and type				
$e^+ e^- \to \tilde{u}_i  \tilde{u}_j$	real(dp) :: SigSup(6,6)				
$e^+  e^- \to \tilde{d}_i  \tilde{d}_j$	real(dp) :: SigSdown(6,6)				
$e^+  e^- \to \tilde{l}_i  \tilde{l}_j$	real(dp) :: SigSle(6,6)				
$e^+ e^- \to \tilde{\nu}_i  \tilde{\nu}_j$	real(dp) :: SigSn(6,6)				
$e^+e^- \to \tilde{\chi}_k^0  \tilde{\chi}_n^0$	real(dp) :: SigChiO(4,4)				
$e^+e^- \to \tilde{\chi}_r^+  \tilde{\chi}_s^-$	real(dp) :: SigN(4,4)				
$e^+e^- \rightarrow h^0 Z, H^0 Z$	real(dp) :: SigSO(2)				
$e^+ e^- \to h^0 A^0, H^0 A^0$	real(dp) :: SigSP(2)				
$e^+  e^-  \rightarrow H^+ H^-$	real(dp) :: SigHp				

```
! output according to SUSY Les Houches Accord
Call LesHouches_Out(HighScaleModel, M_GUT, BRbtosgamma, a_mu, Delta_Rho & & , Ecms, Pm, Pp, ISR, SigSup, SigSdown, SigSle, SigSn, SigChi0 & & , SigC, SigS0, SigSP, SigHp)
! output according to orginal SPheno style
Call WriteOutPutO(11, kont, HighScaleModel, M_GUT, BRbtosgamma, a_mu & & , Delta_Rho, Ecms, Pm, Pp, ISR, SigSup, SigSdown, SigSle , SigSn & & , SigChi0, SigC, SigS0, SigSP, SigHp)
```

The routines SetWriteMinBR and SetWriteMinSig can be used to set a lower on the branching ratios and cross sections [in fb],respectively, which are written to the output files. The routine LesHouches\_Out gives the output according to the SUSY Les Houches Accord [28] whereas the routine WriteOutPutO gives it in the original SPheno style as in Appendix D.

The last statement closes all open files.

call closing() ! closes the files

## 8 Conclusions

We have described SPheno, a program calculating the spectrum, branching ratios and cross sections of supersymmetric particle in  $e^+e^-$  annihilation within the MSSM. The user can choose between the following high scale models: minimal supergravity, minimal supergravity including right handed neutrinos, gauge mediated supersymmetry breaking, anomaly mediated supersymmetry breaking, and string effective field theories based on OI and OII compactification. The calculation of the spectrum are done using two-loop renormalization group equations and the complete one-loop formulas for the SUSY masses. In case of the neutral Higgs bosons and the  $\mu$  parameter leading two-loop effects are included. The masses and mixing angles are used to calculate the most important two body and three body decay modes. They are also used for the calculation of the SUSY production cross sections in  $e^+e^-$  annihilation. Here the effects of initial state radiation and longitudinal beam polarization is included. Finally the following low energy quantities are calculated:  $BR(b \to s\gamma)$ , the supersymmetric contributions to the anomalous magnetic moment of the muon  $a_{\mu}$  and the sfermion contributions to the  $\rho$  parameter. Starting with version 2.2.0 SPheno allows for input and output according to the SUSY Les Houches accord [28].

The program is set up in such a way that extensions can easily be included. The plans for upcoming versions are to include complex phases for the supersymmetric parameters, to include generation mixing, to include QCD and Yukawa corrections for various processes such as Sfermion and Higgs production and decays. In addition beam strahlung for various collider designs will be implemented.

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## A Switches

In this appendix we describe the switches for influencing the behaviour of the program. Inside the main program one can set two scales:

1. The electroweak scale  $M_{EWSB}$ , which is the scale where the loop contributions to the masses and mixing matrices are calculated. The default is to calculate this scale from  $M_{EWSB} = \sqrt{m_{\tilde{t}_1} m_{\tilde{t}_2}}$ . By calling

```
Call SetRGEScale(1.e3_dp**2)
```

 $M_{EWSB}$  will be set to the fixed value of  $10^3$  GeV in this example. Note that the input is the scale squared. In the case one uses a zero or a negative number as input for SetRGEScale then the scale will be calculated from the stop masses.

2. The high energy scale, where the boundary conditions of the model under study are set. Except for GMSB, where the scale is fixed by default, the high scale is calculated from the requirement  $g_1 = g_2$ . By calling

```
Call SetGUTScale(2.e16_dp)
```

 $M_{GUT}$  will be set to  $2 \cdot 10^{16}$  GeV in this example. In the case one uses a zero or a negative number as input for **SetGUTScale** then the scale will be calculated from the requirement  $g_1 = g_2$  except for GMSB.

In general the strong coupling  $g_s$  will be different from  $g_1$  and  $g_2$  in GUT theories if one works at the two loop level [78]. In case someone wants to enforce strict universality at the high scale, this can be done by using the following statement:

```
test = SetStrictUnification(.TRUE.)
```

where test is a logical variable to which the old value of SetStrictUnification is assigned. The same functions can also be used to set this option .FALSE..

Starting with version 2.2.0 it also possible to set these values using the SUSY Les Houches accord as described in Appendix C.

## B Input files

In this section the input files are described. Among these files only the file HighScale.in has to be provided by the user. The other files Control.in, CrossSections.in and StandardModel.in can be used to change the default values which are given below in Appendices B.1, B.2 and B.4. An alternative way to provide SPheno with input is by the SUSY Les Houches accord as described in Appendix C.

## B.1 Control.in

This file contains three entries as shown below:

0 ! ErrorLevel

.True. ! Calculation of branching ratios .True. ! Calculation of cross sections

The values given above are the default values inside the program and are used if the the file Control.in is not present. Here ErrorLevel is an integer in the range [-2,2] where the numbers correspond to the following behaviour of the program:

- -2 do not print severe warnings
- -1 do not print warnings
- 0 print every warning
- 1 abort in case of a severe warning
- 2 abort even in case of a warning

A warning is called severe if either a result is unphysical or if a numerical procedure gives an unreliable result.

## B.2 CrossSections.in

This file contains four entries as shown below:

```
500. ! c.m.s. energy in GeV
```

0. ! degree of longitudinal polarization of electrons0. ! degree of longitudinal polarization of positrons.True. ! calculation of initial state radiation if .TRUE.

The values given above are the default values inside the program and are used if the the file CrossSections.in is not present. Starting with version 2.2.2 up to 100 combinations of the above information will be calculate for a given spectrum. In such a case one has to repeat the above lines as often as required.

## B.3 HighScale.in

In this section we describe the input file for the high scale boundary conditions. The package contains several files starting with Highscale.in and ending in the models described below. One has to rename the model file to the name Highscale.in to use it as input for SPheno. Note that in all examples below the value of  $A_0$  given below will be multiplied by the Yukawas coupling at the high scale and that this product enters the RGEs. For most of the examples below we have used the so-called SPS points defined in [79]. In Appendix D we display the output for the point SPS1a.

#### B.3.1 mSUGRA

The minimal SUGRA version is defined by four parameters and the sign of the  $\mu$  parameter. The parameters are the gaugino mass parameter  $M_{1/2}$ , the scalar mass parameter  $M_0$ , the trilinear parameter  $A_0$  as well as  $\tan \beta$ . The file reads for example

## B.3.2 mSUGRA including right handed neutrinos

In this case one needs four more input values compared to the case of mSUGRA described above: a common right handed neutrino mass  $m_{\nu_R}$  and the light neutrino masses  $m_{\nu_i}$  (i = 1, 2, 3). The file reads for example

```
mSugra
 250.
                                 ! M_1/2
 100.
                                 ! M_0
-100.
                                 ! A_0
                                 ! tan(beta)
  10.
   1.
                                 ! sign of mu
1.e14
                                 ! m_nu_R
1.e-14 3.e-12 0.06e-9
                                 ! m_nu_i
                                 ! if 2-loop RGEs should be used
.TRUE.
```

#### **B.3.3 GMSB**

The GMSB scenario is characterized by two mass parameters  $M_M$  and  $\Lambda$ ; the multiplicity  $N_5$  and  $N_{10}$  of messengers in the  $5+\overline{5}$  and  $10+\overline{10}$  vector-like multiplets, respectively;  $\tan \beta$  and the sign of  $\mu$  as described in Sect. 2.3.3. In addition one can set a common value for the A parameters at the scale  $M_M$ . Note, however, that in the minimal model this value is practically zero. The file reads for example

```
GMSB
100000.
                                   ! Lambda
200000.
                                   ! M_M
1
                                   ! N_5
0
                                   ! N_10
                                   ! AO
0.
                                   ! tan(beta)
15.
                                   ! sign of mu
1.
.TRUE.
                                   ! if 2-loop RGEs should be used
```

#### **B.3.4** AMSB

The implemented AMSB scenario is characterized by the gravitino mass  $m_{3/2}$ , a common scalar mass  $M_0$ ,  $\tan \beta$  and the sign of  $\mu$ . The file reads for example

AMSB ! model 60000. ! M\_3/2 450. ! M\_0 10. ! tan(beta) 1. ! sign of mu .TRUE. ! if 2-loop RGEs should be used

## B.3.5 String I

This scenario is characterized by the gravitino mass  $m_{3/2}$ , the common vev < t > of the moduli fields, the string coupling squared  $g_s^2$ , the sine squared of the mixing angle between the dilaton fields and moduli fields  $\sin^2 \theta$ , the parameter  $\delta_{GS}$  of the Green–Schwarz counter-term, the modular weights  $n_i$  characterizing the couplings between moduli fields and matter fields, which are assumed to be generation independent in the current implementation. Moreover, one needs to specify  $\tan \beta$  and the sign of  $\mu$ . The file reads for example as

```
String_OI
180.
                             ! M_3/2
14.0
                             ! <t>
0.5
                             ! g_s^2
0.9
                             ! sin^2(theta)
 0.
                             ! delta_GS
-1 -3
                             ! n_E n_L
 1 - 2 0
                             ! n_D n_U n_Q
-1 -1
                             ! n_H1 n_H2
10.
                             ! tan(beta)
-1.
                             ! phase(mu)
.TRUE.
                             ! if 2-loop RGEs should be used
```

## B.3.6 String II

This scenario is characterized by the gravitino mass  $m_{3/2}$ , the common vev of the moduli fields < t >, the string coupling squared  $g_s^2$ , the sine squared of the mixing angle between the dilaton fields and moduli fields  $\sin^2 \theta$ , the parameter of the Green–Schwarz counterterm  $\delta_{GS}$ . Moreover, one needs to specify  $\tan \beta$  and the sign of  $\mu$ . There are two different scenarios implemented denoted as String\_OIIa and String\_OIIb corresponding to boundary conditions (A) and (B) of [33], respectively. The files read for example as

```
5.
                             ! tan(beta)
                             ! phase(mu)
1.
.TRUE.
                             ! if 2-loop RGEs should be used
and
String_OIIb
                             ! M_3/2
300.
14.6
                             ! <t>
0.5
                             ! g_s^2
0.9
                             ! sin^2(theta)
                             ! delta_GS
0.
                             ! tan(beta)
5.
1.
                             ! phase(mu)
.TRUE.
                             ! if 2-loop RGEs should be used
```

#### B.3.7 SUGRA

This input file serves as interface for more general models with gauge couplings unification. Here the user can specify non-universal gaugino masses at the high scale  $M_{1/2}[U(1)]$ ,  $M_{1/2}[SU(2)]$ ,  $M_{1/2}[SU(3)]$ , 15 different values of the sfermion mass parameters for every type of sfermions:  $M_{\tilde{E},ii}^0$ ,  $M_{\tilde{L},ii}^0$ ,  $M_{\tilde{D},ii}^0$ ,  $M_{\tilde{U},ii}^0$ ,  $M_{\tilde{Q},ii}^0$ ; two Higgs mass parameters  $M_{H_1}^0$  and  $M_{H_2}^0$ ; nine different A parameters  $A_{0,e,ii}$ ,  $A_{0,d,ii}$  and  $A_{0,u,ii}$ . Here ii denotes that only the diagonal entries can be set.

```
Sugra
480. 300. 300.
                       ! M_1/2_i
150. 150. 150.
                       ! MO_E_ii
150. 150. 150.
                       ! MO_L_ii
150. 150. 150.
                       ! MO_D_ii
150. 150. 150.
                       ! MO_Q_ii
150. 150. 150.
                       ! MO_U_ii
150. 150.
                       ! MO_H_i
0. 0. 0.
                       ! AO_u_ii
0. 0. 0.
                       ! A0_d_ii
0. 0. 0.
                       ! A0_e_ii
                       ! tan(beta)
10.
                       ! phase(mu)
1.
                       ! if 2-loop RGEs should be used
.TRUE.
```

#### **B.3.8** MSSM

This input file serves for the cases that one wants to start with low energy parameters for the calculation of masses, decays and/or production rates. The input consists of the gaugino mass parameters  $M_1$ ,  $M_2$  and  $M_3$ , 15 sfermion mass parameters  $M_{\tilde{E},ii}$ ,  $M_{\tilde{L},ii}$ ,  $M_{\tilde{D},ii}$ ,  $M_{\tilde{U},ii}$ ,  $M_{\tilde{Q},ii}$ , 9 trilinear parameters  $A_{e,ii}$ ,  $A_{d,ii}$ ,  $A_{u,ii}$ ; tan  $\beta$ . Moreover, one has to specify the renormalisation scale Q where all the parameters, including tan  $\beta$ , are given as well as  $\mu$  and the mass of the pseudoscalar Higgs at tree level:  $m_A = B\mu/(\sin\beta\cos\beta)$ .

```
MSSM
99.13 192.74 580.51
                                               ! M_1 M_2 M_3
136.23 136.23 133.55
                                               ! M_E_i
196.64 196.64 195.75
                                               ! M_L_i
519.53 519.53 516.86
                                               ! M_D_i
539.86 539.86 495.91
                                               ! M_Q_i
521.66 521.66 424.83
                                               ! M_U_i
0. 0. -510.01
                                               ! A_u
0.0.-772.66
                                               ! A_d
0. 0. -254.20
                                               ! A e
10. 454.65
                                               ! tan(beta) Q
352.39 393.63
                                               ! mu m_A
```

In the case that the keyword MSSMtree is used instead of MSSM the masses are calculated using tree-level formulas instead of loop corrected masses. The renormalisation scale Q is absent in this case, e.g. the two last line read as

```
10. ! tan(beta) 352.39 393.63 ! mu m_A
```

In addition there is a model file with keyword pMSSM where the following parameters are interpreted differently compared to the case with keyword MSSM:

- $m_A$  is the pole-mass of the pseudoscalar Higgs boson
- $\tan \beta$  is defined to be the value at  $m_Z$

This type of model file has been used to perform the calculations in [75].

## B.4 Standard Model.in

This file contains the values of the Standard Model parameters and must include all lines given below. Otherwise the default values given in the listing below are used:

```
91.1876
                            ! m_Z
2.4952
                            ! Gamma_Z
0.0336 0.0336 0.037
                            ! Br(Z -> 1 1)
0.2
                            ! Br(Z -> invisible)
2.118
                            ! Gamma_W
0.1 0.1 0.1
                            ! Br(W -> 1 nu)
0.51099890e-3
                            ! m_e
0.105658357
                            ! m_mu
1.7770
                            ! m_tau
2.00
                  ! scale Q where the masses of the light quarks u,d,s,c are given
0.003
                  ! m_u(Q)
1.2
                  ! m_c(Q)
174.3
                  ! m_t, pole mass
0.007
                  ! m_d(Q)
                  ! m_s(Q)
0.12
```

```
4.20
                  ! m_b(m_b)
137.0359998
                  ! 1./ alpha
0.1172
                  ! alpha_s(m_Z)
1.16639e-5
                    G_F, Fermi constant
0.224
                 ! s12 of CKM particle data book 1998, 90%, 0.217-0.222
0.0413
                 ! s23 of CKM 0.036-0.042
0.00363
                   ! s13 of CKM 0.0018 - 0.0044
0.
                  ! phase of CKM
                                   0.-2 Pi
2.19709e-6
                  ! life time of muon
3.4e-13
                 ! life time of tau
```

All masses are given in GeV and Br denotes "branching ratio" in the list above.

# C Implementation of SUSY Les Houches Accord

Starting with version 2.2.0 SPheno allows for input and output according to the SUSY Les Houches accord [28]. The name of the input file is LesHouches.in and the output will be written to the file SPheno.spc. In the following we summarize unsupported features as well extensions of this standard. The unsupported features are:

- In block EXTPAR the entries 51–53 are ignored as the corresponding formulas are not (yet) implemented in SPheno.
- Currently there is no information concerning warnings and errors in block SPINFO. This will be changed within the one of the next versions. Please check the file Messages.out for this information.

The current implementation requires that the block MODSEL is read in before the block EXTPAR is read in. In the case that an unknown entry appears, a warning message is printed and SPheno tries to proceed. SPheno stops execution in the case that the model input is not complete.

For the SPheno specific input the block SPhenoInput has to be used. In this block switches can be set, SM input beside the one of the block SMINPUTS can be set. Moreover, the information for the cross section calculation can be given here. Starting with version 2.2.2 several cross sections for different energies and/or polarisation can be calculated for a given spectrum. For this, the entries 22-26 have to be repeated as often as required (there is an upper limit of 100 combinations) and each block has to start with entry 22. Within this block the following flags and parameters can be set, with general structure id value:

```
setting the error level as described Appendix B.1
 2
       if value=1 then the spectrum will be calculated according to the SPA conventions [80]
       if value=1 (0) then (no) branching ratios are calculated
11
12
       only branching ratios larger than value are written out
21
       if value=1 (0) then (no) cross sections are calculated
       cms energy for e^+e^- annihilation, has to be given before entries 23-26
22
       value gives degree of polarisation for e^- beam
23
       value gives degree of polarisation for e^+ beam
24
       if set 1 then ISR corrections will be included, default is 0
25
       only cross sections larger than value (in fb) are written out
26
31
       a fixed value for the GUT scale is used if value is larger than 0
32
       if value=0 then g_3(m_{GUT}) can be different from g_1(m_{GUT}) = g_2(m_{GUT});
       if value=1 then strict unification g_1(m_{GUT}) = g_2(m_{GUT}) = g_3(m_{GUT}) is enforced
       a fixed value for the renormalization scale Q_{EWSB} is used if value is larger than 0
33
41
       sets value of Z-boson width
42
       sets value of W-boson width
   : sets value of electron mass
51
52
   : sets value of muon mass
   : sets scale where the running masses for light quarks (u, d, s, c) are defined
61
62
   : sets value of u-quark mass
63
   : sets value of c-quark mass
   : sets value of d-quark mass
```

# SPheno specific input Block SPhenoInput 1 -1 # error level 11 # calculate branching ratios 12 1.0000000E-04 # write only branching ratios larger than this value 21 # calculate cross section 1 5.0000000E+02 22 # cms energy in GeV 23 0.0000000E+00 # polarisation of incoming e- beam 24 0.0000000E+00 # polarisation of incoming e+ beam 25 # include ISR in the calculation 26 1.0000000E-04 # write only cross sections larger than this value [fb] -1.0000000E+00 31 # m\_GUT, if < 0 than it determined via g\_1=g\_2 32 # require strict unification g\_1=g\_2=g\_3 if '1' is set # Q\_EWSB, if < 0 than Q\_EWSB=sqrt(m\_~t1 m\_~t2)</pre> 33 -1.0000000E+00 # width of the Z-boson 41 2.49520000E+00 42 # width of the W-boson 2.11800000E+00 51 5.10998900E-04 # electron mass 52 1.05658357E-01 # muon mass 61 2.0000000E+00 # scale where quark masses of first 2 gen. are defined 62 3.0000000E-03 # m\_u(Q) 63 1.2000000E+00 # m\_c(Q) 64 7.0000000E-03  $# m_d(Q)$ 

sets value of s-quark mass

65

1.2000000E-01

Here is an example for this block with the default values of SPheno

 $# m_s(Q)$ 

For the output the extensions below have been defined. The information concerning the cross section is written out using a SPheno specific block called SPhenoCrossSections. The first line of this block gives the information on the cms energy, the polarization of the incoming beams as well if ISR is included or not, for example for  $\sqrt{s} = 500 \text{ GeV}$  and unpolarized beams:

```
Block SPhenoCrossSections # cross sections XS 11 -11 500.0 0.00 0.00 1 # e+ e- XS, Pe-, Pe+, including ISR
```

The FORTRAN format is in this case:

```
Format("XS 11 -11 ",F7.1," ",F5.2," ",F5.2," ",A)
```

The cross sections (in fb) themself are written as

Here the first entry gives the cross section in fb, the second entry specifies the number of produced particles, the subsequent two integers give the PDG code of the particles. We have used the FORTRAN format

```
Format(3x, 1P, e16.8, 0p, 3x, I2, 3x, 2(i9, 1x), 2x, " # ", A)
```

As an example we give the cross sections for the SPS1a scenario at an 500 GeV  $e^+e^-$  linear collider with unpolarized beams:

```
Block SPhenoCrossSections
                            # cross sections
XS 11 -11
             500.0 0.00
                          0.00 1
                                   # e+ e- XS, Pe-, Pe+,
                                                             including ISR
#
                         NDA
       Sigma [fb]
                                     ID1
                                              ID2
                                                     # ~e_R-
     2.83574498E+02
                        2
                                        -2000011
                                                                    ~e_R+
                               2000011
                        2
                                                     # ~e_R-
                                                                    ~e_L+
     7.79728001E+01
                               2000011
                                        -1000011
     4.57495061E+01
                        2
                               1000011
                                        -1000011
                                                     # ~e_L-
                                                                    ~e_L+
                                                                    ~mu_R+
     5.47916441E+01
                        2
                               2000013
                                        -2000013
                                                     # ~mu_R-
     6.00045490E-03
                        2
                               2000013
                                        -1000013
                                                       ~mu_R-
                                                                    ~mu_L+
     1.90114309E+01
                        2
                               1000013
                                        -1000013
                                                     # ~mu_L-
                                                                    ~mu_L+
                        2
                                                     # ~tau_1-
     5.96228076E+01
                               1000015
                                        -1000015
                                                                    ~tau_1+
                        2
     1.26426385E+00
                               1000015
                                        -2000015
                                                     # ~tau_1-
                                                                    ~tau_2+
                        2
     1.59684572E+01
                               2000015
                                        -2000015
                                                     # ~tau_2-
                                                                    ~tau_2+
                        2
                                                     # ~nu_eL
     4.52889205E+02
                               1000012
                                        -1000012
                                                                    ~nu_eL*
     1.36168303E+01
                        2
                               1000014
                                        -1000014
                                                     # ~nu_muL
                                                                    ~nu_muL*
                        2
     1.39168830E+01
                               1000016
                                        -1000016
                                                     # ~nu_tauL
                                                                    ~nu_tauL*
     2.75869582E+02
                        2
                               1000022
                                          1000022
                                                     # chi_10 chi_10
                        2
     6.56937491E+01
                               1000022
                                          1000023
                                                     # chi_10 chi_20
                        2
                                                     # chi_10 chi_30
     7.10141133E+00
                               1000022
                                          1000025
                        2
                               1000022
                                                     # chi_10 chi_40
     8.27993814E-01
                                          1000035
                        2
     6.90281358E+01
                               1000023
                                          1000023
                                                     # chi_20 chi_20
     1.60903760E+02
                        2
                               1000024
                                         -1000024
                                                     # chi_1- chi_1+
     2.47077869E+01
                        2
                                    25
                                               23
                                                     # h0 Z
```

The information concerning the value of low energy observables  $(BR(b \to s\gamma), SUSY)$  contribution to  $(g-2)_{\mu}$  and  $\Delta(\rho)$  is written out using a SPheno specific block called SPhenoLowEnergy. We use the following identifier:

1 : BR( $b \rightarrow s\gamma$ )

2 : SUSY contributions to  $(g-2)_{\mu}$ 

3 : SUSY contributions to  $\Delta(\rho)$ 

As an example we give here the output for the SPS1a scenario:

Block SPhenoLowEnergy # low energy observables

- 1 4.55809155E+00 # BR(b -> s gamma)
- 2 5.42193822E-09 # (g-2)\_muon
- 3 1.97608480E-04 # Delta(rho)

## D Sample output

Here we give the content of the file SPheno.out provided one uses the content of HighScale.in for the mSUGRA scenario described in Appendix B.3.1 and the default values of the files Control.in, CrossSections.in and StandardModel.in.

SPheno output file

Version 2.2.2, created: 14.09.2004, 17:14

mSugra input at the GUT scale 2.4620574378552756E+16

M\_1/2 : 2.5000000000000000E+02
M\_0 : 1.00000000000000E+02
A\_0 : -1.000000000000000E+02
tan(beta) at m\_Z : 10.000000000000000

Running masses have been used for the boundary conditions at mZ

Parameters at the scale 4.8442121445544416E+02

g' g g\_3

3.61098068E-01 6.46530088E-01 1.09487945E+00

Y\_e Y\_mu Y\_tau

2.88425398E-05 5.96372335E-03 1.00306479E-01

 $Y_u$   $Y_c$   $Y_t$ 

8.84588597E-06 3.53835431E-03 8.92075460E-01

Yd Ys Yb

1.91353683E-04 3.28034867E-03 1.37341541E-01

Gaugino mass parameters

1.0155024266722937E+02 1.9167894566112778E+02 5.8533134393167563E+02

```
mu, B
  3.5775913791766874E+02 1.6731090082362174E+04
Slepton mass parameters
A_1
-2.5344637795775390E+02 -2.5344027822464955E+02 -2.5172098902647821E+02
M2_E
  1.8443988295117768E+04 1.8441668307344480E+04 1.7789728306764348E+04
M2 L
 3.8222680589380434E+04
                          3.8221538628557908E+04
                                                   3.7900699251843856E+04
Squark mass parameters
A_d
-8.5558462724494541E+02 -8.5558116731465941E+02 -7.9140619867408907E+02
 -6.8014557244910873E+02 -6.8014185758548206E+02 -4.9696697403034966E+02
M2_D
  2.7491167922793247E+05 2.7490965928240586E+05
                                                   2.7157606836767372E+05
M2_U
  2.7715958460406726E+05 2.7715762927385850E+05
                                                   1.7637959938821895E+05
M2 Q
  2.9720857969470561E+05 2.9720661515585968E+05
                                                   2.4639278946889582E+05
Higgs mass parameters
  3.2551122776108325E+04 -1.2811756834297138E+05
Masses and mixing matrices
Gluino :
          6.0453116356737542E+02
                                    1.00000000000000000
Charginos
  1.8029681792857122E+02 3.8336666782688877E+02
       -0.91584
                        0.40153
                        0.91584
        0.40153
  V
        -0.97278
                        0.23175
        0.23175
                        0.97278
Neutralinos
 97.0684422442318606
                      1.8069642547695042E+02 3.6506776000040759E+02
                                                                       3.82276404251164
  N
   1
      1 (
                  -0.98582,
                                    0.00000)
   1
     2 (
                  0.05596,
                                    0.00000)
      3 (
                  -0.14856,
                                    0.00000)
   1
     4 (
                   0.05430,
                                    0.00000)
                  -0.10355,
                                    0.00000)
```

2	2	(	-0.94298,	0.00000)
2	3	(	0.27441,	0.00000)
2	4	(	-0.15735,	0.00000)
3	1	(	0.00000,	0.06043)
3	2	(	-0.00000,	-0.09021)
3	3	(	-0.00000,	-0.69486)
3	4	(	-0.00000,	-0.71090)
4	1	(	0.11737,	0.00000)
4	2	(	-0.31546,	0.00000)
4	3	(	-0.64792,	0.00000)
4	4	(	0.68331,	0.00000)

e-sneutrino mass : 1.8629967789333597E+02 mu-sneutrino mass : 1.8629643037340736E+02 tau-sneutrino mass : 1.8538056982341820E+02

#### selectron masses

1.4394583534020416E+02 2.0249814925123943E+02

R\_e

0.00009 1.00000 -1.00000 0.00009

#### smuon masses

1.4391005493518603E+02 2.0251437533591726E+02

 $R_{mu}$ 

0.01958 0.99981 -0.99981 0.01958

#### stau masses

1.3430774918769018E+02 2.0648286202975819E+02

R\_tau

0.28332 0.95903 -0.95903 0.28332

## u-squark masses

5.4818625728877964E+02 5.6590020431737230E+02

 $R_u$ 

0.00006 1.00000 -1.00000 0.00006

## c-squark masses

5.4817448074068113E+02 5.6590853096426122E+02

 $R_c$ 

0.02385 0.99972 -0.99972 0.02385

## t-squark masses

3.9989424629928800E+02 5.8682205427167264E+02

R\_t

0.55322 0.83304 -0.83304 0.55322

d-squark masses

5.4791161728759471E+02 5.7129911929796879E+02

R\_d

0.00058 1.00000 -1.00000 0.00058

s-squark masses

5.4790724305571359E+02 5.7129944562180071E+02

 $R_s$ 

0.00999 0.99995 -0.99995 0.00999

b-squark masses

5.1564470876393955E+02 5.4769779167739352E+02

 $R_b$ 

0.94719 0.32066 -0.32066 0.94719

 $m_AO$ ,  $m_H+$ 

3.9982837260329677E+02 4.0811814345802685E+02

m\_h0, m\_H0

1.1082357413611736E+02 4.0020427017852177E+02

R\_S0

0.11369 0.99352 -0.99352 0.11369

Low energy constraints

10^4 Br(b -> s gamma) : 0.4581123E+01 Delta(a\_mu) : 0.5696755E-08 Delta(rho) : 0.2001459E-03

Anti particles are marked with a \* in case of (s)neutrinos and (s)quarks in the decay section.

Decay widths (GeV) and branching ratios

Selectron\_1

Neutralino\_1 e 0.21291502 100.00000000

Total width : 0.21291502

Selectron_2 Neutralino_1 e Neutralino_2 e Chargino_1 neutrino Total width :	0.12098066 0.03480167 0.06033029 0.21611263	
Smuon_1 Neutralino_1 mu Total width:	0.21263340 0.21263340	100.00000000
Smuon_2 Neutralino_1 mu Neutralino_2 mu Chargino_1 neutrino Total width :	0.12115824 0.03482867 0.06037630 0.21636321	16.09731436
Stau_1 Neutralino_1 tau Total width :	0.14502344 0.14502344	100.00000000
Stau_2 Neutralino_1 tau Neutralino_2 tau Chargino_1 neutrino Total width :	0.15977947 0.04134796 0.07135516 0.27248259	15.17453244
e-Sneutrino Neutralino_1 neutrino Neutralino_2 neutrino Chargino_1 e Total width :	0.14973207 0.00211800 0.00585751 0.15770758	94.94284818 1.34299390 3.71415792
<pre>mu-Sneutrino Neutralino_1 neutrino Neutralino_2 neutrino Chargino_1 mu Total width :</pre>	0.14972554 0.00211562 0.00584974 0.15769090	

tau-Sneutrino

Neutralino_1 neutrino Neutralino_2 neutrino Chargino_1 tau Total width :	0.14788387 0.00149472 0.00380717 0.15318576	96.53891386 0.97575508 2.48533106
Sdown_1 Neutralino_1 d-quark Neutralino_2 d-quark Neutralino_3 d-quark Neutralino_4 d-quark Total width :	0.28838606 0.00269444 0.00035714 0.00114758 0.29258676	98.56429027 0.92090214 0.12206173 0.39221887
Sdown_2 Neutralino_1 d-quark Neutralino_2 d-quark Neutralino_3 d-quark Neutralino_4 d-quark Chargino_1 u-quark Chargino_2 u-quark Total width:	0.12870838 1.64327451 0.00857068 0.08252004 3.23407423 0.23170134 5.32884919	2.41531292 30.83732443 0.16083549 1.54855280 60.68991849 4.34805586
S-strange_1 Neutralino_1 s-quark Neutralino_2 s-quark Neutralino_3 s-quark Neutralino_4 s-quark Chargino_1 c-quark Total width:	0.28839626 0.00291737 0.00036637 0.00114598 0.00045256 0.29327899	98.33512617 0.99474185 0.12492215 0.39074858 0.15431122
S-strange_2 Neutralino_1 s-quark Neutralino_2 s-quark Neutralino_3 s-quark Neutralino_4 s-quark Chargino_1 c-quark Chargino_2 c-quark Total width:	0.12869979 1.64305243 0.00859993 0.08255413 3.23357116 0.23179659 5.32827402	2.41541244 30.83648512 0.16140172 1.54935962 60.68702822 4.35031288
Sbottom_1 Neutralino_1 b-quark Neutralino_2 b-quark Neutralino_3 b-quark	0.16653332 1.34625440 0.01956140	4.31243760 34.86172159 0.50654921

Neutralino_4 b-quark Chargino_1 t-quark Stop_1 W- Total width :	0.04233574 1.72259277 0.56442035 3.86169797	1.09629870 44.60713339 14.61585951
Sbottom_2 Neutralino_1 b-quark Neutralino_2 b-quark Neutralino_3 b-quark Neutralino_4 b-quark Chargino_1 t-quark Stop_1 W- Total width:	0.24058254 0.09329885 0.04185389 0.05807690 0.12191203 0.19798439 0.75370860	31.91983477 12.37863659 5.55305967 7.70548478 16.17495579 26.26802841
Sup_1 Neutralino_1 u-quark Neutralino_2 u-quark Neutralino_3 u-quark Neutralino_4 u-quark Total width:	1.15420363 0.01078281 0.00143143 0.00460140 1.17101929	98.56401566 0.92080534 0.12223802 0.39293982
Sup_2 Neutralino_1 u-quark Neutralino_2 u-quark Neutralino_3 u-quark Neutralino_4 u-quark Chargino_1 d-quark Chargino_2 d-quark Total width:	0.03614750 1.75889152 0.00500542 0.06001133 3.59863170 0.07406777 5.53275523	0.65333627 31.79051757 0.09046880 1.08465542 65.04230795 1.33871398
S-charm_1 Neutralino_1 c-quark Neutralino_2 c-quark Neutralino_3 c-quark Neutralino_4 c-quark Chargino_1 s-quark Total width :	1.15355953 0.01183744 0.00147948 0.00458567 0.00217222 1.17363442	98.28951064 1.00861420 0.12605944 0.39072421 0.18508491
S-charm_2 Neutralino_1 c-quark Neutralino_2 c-quark Neutralino_3 c-quark	0.03677444 1.75780413 0.00500431	0.66496763 31.78519855 0.09048962

Neutralino_4 c-quark Chargino_1 s-quark Chargino_2 s-quark Total width :	0.06008384 3.59645144 0.07414186 5.53026002	1.08645601 65.03223042 1.34065776
Stop_1 Neutralino_1 t-quark Neutralino_2 t-quark Chargino_1 b-quark Chargino_2 b-quark c-quark neutralino_1 c-quark neutralino_2 Total width:	0.39786160 0.24313066 1.37082175 0.01991200 0.00040360 0.01643838 2.04857653	19.42136860 11.86827336 66.91581865 0.97199197 0.01970164 0.80242915
Stop_2 Neutralino_1 t-quark Neutralino_2 t-quark Neutralino_3 t-quark Neutralino_4 t-quark Chargino_1 b-quark Chargino_2 b-quark Stop_1 Z Stop_1 h0 Total width:	0.22193261 0.64232737 0.30940040 1.44196179 1.62955414 1.44485203 1.39455443 0.26950476 7.35408754	3.01781302 8.73429056 4.20718955 19.60762341 22.15848167 19.64692457 18.96298387 3.66469336
Chargino_1 Smuon_1 neutrino Stau_1 neutrino Neutralino_1 W neutralino_1 e^+ nu neutralino_1 mu^+ nu neutralino_1 tau^+ nu Total width :	0.00004613 0.01510357 0.00071260 0.00003745 0.00003745 0.00003773 0.01597497	0.28874351 94.54518351 4.46072016 0.23441594 0.23441507 0.23615356
Chargino_2 Selectron_2 neutrino Smuon_2 neutrino Stau_1 neutrino Stau_2 neutrino e-sneutrino e mu-sneutrino mu tau-sneutrino tau Neutralino_1 W Neutralino_2 W	0.13339597 0.13343849 0.00064169 0.14518928 0.04987592 0.04994101 0.06840093 0.16988544 0.73896629	5.23062736 5.23229482 0.02516131 5.69305830 1.95569876 1.95825114 2.68208833 6.66142664 28.97581836

Chargino_1 Z Chargino_1 h0 neutralino_1 b^* t neutralino_2 b^* t chargino_1 b b^* Total width :	0.60911313 0.45051419 0.00030103 0.00003051 0.00054500 2.55028619	23.88410873 17.66524071 0.01180365 0.00119636 0.02137000
Neutralino_1 : stable		
Neutralino_2		
Selectron^1 e^+	0.00066378	3.07448383
Selectron^+_1 e^-	0.00066378	3.07448383
Smuon^1 mu^+	0.00069076	3.19948480
Smuon^+_1 mu^-	0.00069076	3.19948480
Stau^1 tau^+	0.00939369	43.50975897
Stau^+_1 tau^-	0.00939369	43.50975897
neutralino_1 u u^*	0.00000345	0.01599688
neutralino_1 c c^*	0.00000345	0.01597295
neutralino_1 d d^*	0.00000459	0.02123795
neutralino_1 s s^*	0.00000459	0.02123789
<pre>neutralino_1 b b^* neutralino_1 nu_e nu_e^</pre>	0.00000472 0.00001768	0.02186206 0.08190096
<pre>neutralino_1 nu_e    nu_e^ neutralino_1 nu_mu    nu_mu</pre>	0.00001768	0.08190090
neutralino_1 nu_mu nu_mu neutralino_1 nu_tau nu_ta	0.00001769	0.08192737
neutralino_1 e^- e^+	0.00001544	0.02919519
neutralino_1 mu^- mu^+	0.00000630	0.02917454
neutralino_1 tau^- tau^+	0.00000514	0.02379866
Total width :	0.02158986	
Neutralino_3		
Selectron^1 e^+	0.00245868	0.12468320
Selectron^+_1 e^-	0.00245868	
Selectron^2 e^+	0.00115857	
Selectron^+_2 e^-	0.00115857	0.05875282
Smuon^1 mu^+	0.00247970	0.12574917
Smuon^+_1 mu^-	0.00247970	0.12574917
Smuon^2 mu^+	0.00120501	0.06110808
Smuon^+_2 mu^-	0.00120501	0.06110808
Stau^1 tau^+	0.00993392	0.50376391
Stau^+_1 tau^-	0.00993392	0.50376391
Stau^2 tau^+	0.01297595	
Stau^+_2 tau^-	0.01297595	
e-sneutrino nu_e^*	0.00636550	0.32280417
e-sneutrino^* nu_e	0.00636550	0.32280417
mu-sneutrino nu_mu^*	0.00636566	
mu-sneutrino^* nu_mu	0.00636566	0.32281210

tau-sneutrino nu_tau^*	0.00640970	0.32504561
tau-sneutrino^* nu_tau	0.00640970	0.32504561
Chargino^+_1 W^-	0.58233645	29.53114356
Chargino^1 W^+	0.58233645	29.53114356
Neutralino_1 Z	0.22153690	11.23446435
Neutralino_2 Z	0.42050000	21.32417766
Neutralino_1 h0	0.04208071	2.13397535
Neutralino_2 h0	0.02440336	1.23753049
Neutralino_2 photon	0.00002030	0.00102937
Total width :	1.97194005	
N+1 4		
Neutralino_4	0 01002407	0.27554000
Selectron^1 e^+	0.01003487	0.37554208
Selectron^+_1 e^-	0.01003487	0.37554208
Selectron^2 e^+	0.02565497	
Selectron^+_2 e^-	0.02565497	
Smuon^1 mu^+	0.01001475	
Smuon^+_1 mu^-	0.01001475	0.37478909
Smuon^2 mu^+	0.02572432	0.96269931
Smuon^+_2 mu^-	0.02572432	
Stau^1 tau^+	0.00708230	0.26504573
Stau^+_1 tau^-	0.00708230	0.26504573
Stau^2 tau^+	0.04322619	1.61768406
Stau^+_2 tau^-	0.04322619	1.61768406
e-sneutrino nu_e^*	0.06696143	2.50594432
e-sneutrino^* nu_e	0.06696143	2.50594432
<pre>mu-sneutrino nu_mu^*</pre>	0.06696288	2.50599874
mu-sneutrino^* nu_mu	0.06696288	2.50599874
tau-sneutrino nu_tau^*	0.06737263	2.52133329
tau-sneutrino^* nu_tau	0.06737263	2.52133329
Chargino^+_1 W^-	0.68286763	25.55543352
Chargino^1 W^+	0.68286763	25.55543352
Neutralino_1 Z	0.05548528	2.07646431
Neutralino_2 Z	0.05010799	1.87522654
Neutralino_1 h0	0.18248851	6.82939545
Neutralino_2 h0	0.37208545	13.92481435
Neutralino_2 photon	0.00003496	0.00130835
Total width :	2.67210349	
Gluino		
Sup_1 u^*	0.22549273	4.93757980
Sup_1^* u	0.22549273	4.93757980
Sup_2 u^*	0.10927993	2.39288589
Sup_2^* u	0.10927993	2.39288589
S-charm_1 c^*	0.22529588	4.93326946

```
S-charm_1^* c
                                               4.93326946
                                0.22529588
  S-charm_2 c^*
                                0.10935220
                                               2.39446832
  S-charm_2^* c
                                0.10935220
                                               2.39446832
  Stop_1 t^*
                                0.23631434
                                               5.17453901
  Stop_1^* t
                                0.23631434
                                               5.17453901
  Sdown_1 d^*
                                0.22758779
                                               4.98345500
  Sdown_1^* d
                                0.22758779
                                               4.98345500
  Sdown_2 d^*
                                0.08161707
                                               1.78715654
  Sdown_2^* d
                                0.08161707
                                               1.78715654
  S-strange_1 s^*
                                0.22761069
                                               4.98395651
  S-strange_1^* s
                                0.22761069
                                               4.98395651
  S-strange_2 s^*
                                0.08162105
                                               1.78724357
  S-strange_2^* s
                                0.08162105
                                               1.78724357
  Sbottom_1 b^*
                                0.51303196
                                              11.23378233
  Sbottom_1^* b
                                0.51303196
                                              11.23378233
                                0.23942646
                                               5.24268461
  Sbottom_2 b^*
  Sbottom_2^* b
                                0.23942646
                                               5.24268461
  Stop_1 c^*
                                0.00526030
                                               0.11518395
  Stop_1^* c
                                0.00526030
                                               0.11518395
  neutralino_2 gluon
                                0.00021864
                                               0.00478751
  neutralino_3 gluon
                                0.00032667
                                               0.00715313
  neutralino_4 gluon
                                0.00037459
                                               0.00820239
  neutralino_1 t t^*
                                0.00008996
                                               0.00196988
  neutralino_2 t t^*
                                0.00009543
                                               0.00208969
  chargino^+_1 t^* b
                                0.00067696
                                               0.01482320
  chargino^-_1 t b^*
                                0.00067696
                                               0.01482320
  chargino^+_2 t^* b
                                0.00030059
                                               0.00658192
  chargino^-_2 t b^*
                                0.00030059
                                               0.00658192
 Total width:
                                4.56686756
h0
                                0.00000104
                                               0.03717403
  muons
  taus
                                0.00029367
                                              10.49412940
                                0.00000127
                                               0.04524099
  s-quark
                                0.00216519
                                              77.37202572
  b-quark
  c-quark
                                0.00010749
                                               3.84124361
  W+ W-*
                                0.00005707
                                               2.03933591
  W+* W-
                                0.00005707
                                               2.03933591
  Z Z*
                                0.00000615
                                               0.21967100
                                0.00010946
                                               3.91166457
  g g
 Total width:
                                0.00279842
HO
                                0.00028046
                                               0.03480344
  muons
  taus
                                0.07932609
                                               9.84378583
  s-quark
                                0.00031273
                                               0.03880696
```

b-quark t-quark Selectron 1 1 Smuon 1 1 Smuon 1 2 Smuon 2 1 Stau 1 1 Stau 1 2 Stau 2 1 e-Sneutrino mu-Sneutrino neutralino_1 neutralino_1 neutralino_1 neutralino_2 neutralino_2 neutralino_2 chargino^+_1 chargino^1 Z Z W+ W- hO hO g g Total width :	0.54671657 0.03547964 0.00038368 0.00039169 0.00001977 0.00401977 0.00461331 0.00404359 0.00404359 0.00089263 0.00089273 0.00092078 0.01634146 0.04742007 0.01339966 0.03593261 0.00145026 0.00310488 0.00951582 0.00034176 0.80584938	4.40276284 0.04761235 0.04860540 0.00245370 0.00245370 0.57247839 0.50178033 0.50178033 0.11076927 0.11078183 0.11426177 2.02785540 5.88448299 1.66280004 4.45897405 0.17996641 0.38529262 1.18084312
muons taus s-quark b-quark t-quark Smuon 1 2 Smuon 2 1 Stau 1 2 Stau 2 1 neutralino_1 neutralino_1 neutralino_2 neutralino_2 chargino^+_1 chargino^1 h0 Z Total width :	0.00028106 0.07950191 0.00031341 0.54815429 0.11221779 0.00002049 0.00594160 0.00594160 0.02505516 0.10779205 0.09278079 0.24090139 0.00267484 1.22160158	6.50800650 0.02565593 44.87177335
H^+ muon neutrino tau neutrino s-quark c-quark	0.00028679 0.08112135 0.00026690	0.04230494 11.96655479 0.03937192

b-quark t-quark	0.43329892	63.91775887
Selectron_2 Sneutrino	0.00073931	0.10905855
Smuon_1 Sneutrino	0.00005822	0.00858803
Smuon_2 Sneutrino	0.00073338	0.10818334
Stau_1 Sneutrino	0.01538016	2.26879229
<pre>chargino_1 neutralino_1</pre>	0.14190050	20.93234365
<pre>chargino_1 neutralino_2</pre>	0.00107290	0.15826853
hO W	0.00303784	0.44812476
Total width :	0.67790067	

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