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SPheno, a program for calculating supersymmetric spectra, SUSY particle decays and SUSY particle production at e^+e^- colliders $\stackrel{\diamond}{\sim}$

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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 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 μ 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 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.

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PROGRAM SUMMARY	vironment, has also been tested on an HP machine and an ALPHA machine		
Title of program: SPheno	Programming language used: F90		
Catalogue identifier: ADRV	Number of bytes in distributed program, including test data, etc.:		
Program Summary URL: http://cpc.cs.qub.ac.uk/summaries/ADRV	450 615		
Program obtainable from: CPC Program Library, Queen's Univer-	No. of lines in combined program and test case: 43319		
sity of Belfast, N. Ireland	Distribution format: tar gzip file		
Computer for which the program is designed and others on which it is operable: PC running under Linux, should run in every Unix en-	Separate documentation available: no		

* This program can be downloaded from the CPC Program Library under catalogue identifier: http://cpc.cs.qub.ac.uk/summaries/ADRV *E-mail address:* porod@physik.unizh.ch (W. Porod).

0010-4655/03/\$ - see front matter © 2003 Elsevier Science B.V. All rights reserved. doi:10.1016/S0010-4655(03)00222-4 *Keywords:* Supersymmetric Standard Model, renormalization group equations, MSSM spectra, Runge–Kutta, decays of super-symmetric particles, production

Nature of physical problem

The first issue is the determination of the masses and couplings of supersymmetric particles in the R-parity conserved MSSM. Low energy data on Standard Model fermion masses, gauge couplings and electroweak gauge boson masses serve as constraints. Radiative corrections from supersymmetric particles to these inputs must be calculated. Theoretical constraints on the soft SUSY breaking parameters from a high scale theory are imposed and the parameters at the electroweak scale are obtained from the high scale parameters by evaluating the corresponding renormalization group equations. These parameters must be consistent with the requirement of correct electroweak symmetry breaking. The second issue is to use the obtained masses and couplings for calculating decay widths and branching ratios of supersymmetric particles as well as the cross sections for theses particles in electron positron annihilation. The third issue is to calculate the following low energy constraints: the branching ratio for $b \rightarrow s\gamma$, the SUSY contributions to the ρ parameter as well as the SUSY contributions to the anomalous magnetic moment of the muon.

Method of solution

The renormalization connecting a high scale and the electroweak scale is calculated by the Runge–Kutta method. Iteration provides a solution consistent with the multi-boundary conditions. In case of three-body decays and for the calculation of initial state radiation Gaussian quadrature is used for the numerical solution of the integrals.

Restrictions on the complexity of the problem

MSSM spectra with real parameters only neglecting the effect of mixing between (s)fermion generations.

Typical running time: 0.3 s on a Pentium III

LONG WRITE-UP

1. Introduction

Supersymmetry (SUSY) [1–3] 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 [8–11].

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 SPheno¹ 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

¹ SPheno stands for Supersymmetric Phenomenology.

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 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 $b \rightarrow s\gamma$, the supersymmetric contributions to the anomalous magnetic of the muon a_{μ} as well as supersymmetric contributions to the ρ parameter.

For the calculation of the spectrum the programs ISASUSY [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, ISASUSY [17], SPYTHIA [21] and SUSYGEN [22]. A comparison of the results of these programs will be given in a future paper.

SPheno has been written in Fortran90. 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) [23], complete one-loop correction to all SUSY and Higgs masses [24] supplemented by the 2-loop corrections to the neutral Higgs bosons [25,26] and to the μ parameter [26]. The present version of SPheno does all calculations for real parameters neglecting the flavor 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 Section 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 Section 3 we list the implemented decay modes of supersymmetric particles and the Higgs bosons. We also discuss shortly the approximations used. In Section 4 we present the implemented cross sections in $e^+e^$ annihilation. In Section 5 we discuss give details on the implemented low energy constraints. In Section 6 we discuss the implemented algorithm in some detail. In Section 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. Moreover, we list the output of the program for a typical example. The source code as well as precompiled versions of the program can be obtained from the author via email: porod@physik.unizh.ch; or it can be downloaded from http://www-theorie.physik.unizh.ch/~porod/SPheno.html.

2. MSSM parameters, particle spectrum, and high scale 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:

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

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; ϵ_{ab} is the totally antisymmetric tensor with $\epsilon_{12} = 1$. Note that the sign of μ is identical to the one in ISASUSY [17] and SOFTSUSY [18] but opposite to the convention in [24]. 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

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

mass terms for scalar matter fields and Higgs fields

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

and trilinear couplings of scalar matter fields and Higgs fields

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

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

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

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

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

$$X = \begin{pmatrix} M & \frac{g}{\sqrt{2}}v_2\\ \frac{g}{\sqrt{2}}v_1 & \mu \end{pmatrix}.$$
(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 $\tilde{h}_1^0, \tilde{h}_2^0$ form the neutralinos. In the basis $\tilde{\psi}^0 = (\tilde{b}, \tilde{w}^3, \tilde{h}_1^0, \tilde{h}_2^0)^{\mathrm{T}}$ one finds the mass term $-(\tilde{\psi}^0)^{\mathrm{T}} Y \tilde{\psi}^0$ 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 α 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:

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

The other sfermion mass matrices are 2×2 matrices:

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

$$m_{\tilde{f}}^2 = R_{\tilde{f}}^{\dagger} M_{\tilde{f}}^2 R_{\tilde{f}}.$$
(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:

$$|\mu|^{2} = \frac{1}{2} \left[\tan 2\beta \left(M_{H_{2}}^{2} \tan \beta - M_{H_{1}}^{2} \cot \beta \right) - m_{Z}^{2} \right].$$
(18)

Moreover, in all cases the high scale parameters are supplemented by the sign of μ and tan β .

2.3.1. Minimal supergravity

The minimal supergravity (mSUGRA) scenario is characterized by a set of universal parameters [2,13] 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 :

$$M_i(M_{\rm GUT}) = M_{1/2},$$
(19)
$$M_i^2(M_{\rm GUT}) = M_i^2$$
(20)

$$M_{\tilde{j}}^{2}(M_{\rm GUT}) = M_{0}^{2},$$
(20)
 $A_{i}(M_{\rm GUT}) = A_{0}Y_{i}(M_{\rm GUT}).$
(21)

$$m(m001) = m01/(m001)$$
.

2.3.2. Minimal supergravity including right handed neutrinos

In addition to the parameters of the mSUGRA model above the following parameters appear in this case: m_{ν_R} , a common mass for all right handed neutrinos, and m_{ν_i} (i = 1, 2, 3), the light neutrino masses. In this case the MSSM RGEs are run up to the scale m_{ν_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_{ν_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:

$$M_{\tilde{R}}^{2}(M_{\rm GUT}) = M_{0}^{2}, \tag{22}$$

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

The corresponding RGEs are used in the running from M_{GUT} to m_{ν_R} . At the scale m_{ν_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_{ν_R} the usual set of MSSM RGEs are used.

2.3.3. Gauge mediated supersymmetry breaking

Gauge mediated supersymmetry breaking [14,27] (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 [28] 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:

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

with $k_i = 1, 1, 3/5$ for SU(3), SU(2), and U(1), respectively; the coefficients C_j^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 [28]

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

$$M_a = \frac{1}{g_a} \beta_a m_{3/2},$$
(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 β_a and β_{Y_i} are the beta functions of gauge and Yukawa couplings, respectively. The γ_j are the anomalous dimensions of the corresponding matter superfield and $m_{3/2}$ is the gravitino mass. Eq. (30) leads to negative mass squared for the sleptons which is phenomenologically not acceptable. There are several possibilities to solve this problem [29] 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 [30,31].

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 [30,31]. 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 [31]. 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 [31] as well as an OI compactification scheme. For the implementation of the OI compactification scheme we have used formulas (3.21)–(3.23) of [31]:

$$M_{i} = -g_{i}^{2}m_{3/2} \left\{ s\sqrt{3}\sin\theta + \left[b_{i} + s\sqrt{3}\sin\theta g_{s}^{2} \left(C_{i} - \sum_{j} C_{i}^{j} \right) + 2t\cos\theta G_{2}(t) \left(\delta_{\text{GS}} + b_{i} - 2\sum_{j} C_{i}^{j}(1+n_{j}) \right) \right] / 16\pi^{2} \right\},$$
(32)
$$M_{\tilde{j}}^{2} = m_{3/2}^{2} \left\{ (1+n_{j}\cos^{2}\theta) + 2\sqrt{3}s\sin\theta \left[\sum_{j} \gamma_{j}^{i}g_{i}^{2} - \frac{1}{2s}\sum_{j} \gamma_{j}^{km} \right] \right\}$$

$$= m_{3/2} \left\{ (1 + n_j \cos \theta) + 2\sqrt{3} \sin \theta \left[\sum_i \gamma_j g_i - \frac{1}{2s} \sum_{km} \gamma_j \right] + \gamma_j + 2t \cos \theta G_2(t) \sum_{km} \gamma_j^{km} (n_j + n_k + n_m + 3) \right\},$$
(33)

$$A_{jkm} = m_{3/2} \Big[-\sqrt{3}\sin\theta - 2t\cos\theta (n_j + n_k + n_m + 3)G_2(t) + \gamma_j + \gamma_k + \gamma_m \Big],$$
(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 ζ denoting the Riemann zeta function. δ_{GS} is the parameter of the Green–Schwarz counterterm. γ_j are the anomalous dimensions of the

matter fields, the γ_j^i and γ_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:

$$A_{e,n}(\text{GUT}) = Y_{e,nn}(\text{GUT})A_{\tilde{E}_n\tilde{L}_nH_1},$$
(35)

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

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

where n denotes the generation.

In case of the OII compactification scheme the gaugino masses are given by Eqs. (3.11) of [31]:

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

For the sfermion parameters we have implemented two sets of boundary conditions: set (A) is specified by formulas (3.15) and (3.19) of [31]:

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

$$A_{ijk} = m_{3/2} \left\{ \gamma_i + \gamma_j + \gamma_k - \frac{\sqrt{3}k_s \sin\theta}{k_{s\bar{s}}^{1/2}} \right\}.$$
 (40)

Set (*B*) is specified by formulas (3.16) and (3.20) of [31]:

$$M_{i}^{2} = m_{3/2}^{2} \left\{ \frac{\sqrt{3} \sin \theta}{k_{s\bar{s}}^{1/2}} [1 + 2t \cos \theta G_{2}] \left[\sum_{a} g_{a}^{2} \gamma_{i}^{a} + \frac{1}{2} \sum_{jk} \gamma_{i}^{jk} (k_{s} + k_{\bar{s}}) \right] \\ + \sin^{2} \theta \left[1 + \gamma_{i} + \ln[2t|\eta(t)|^{4}] \left(\sum_{a} \gamma_{i}^{a} + 2 \sum_{jk} \gamma_{i}^{jk} \right) - \sum_{a} \gamma_{i}^{a} \ln(g_{a}^{2}) \right] \\ - \frac{9 \sin^{2} \theta}{k_{s\bar{s}}} \left[\sum_{a} \gamma_{i}^{a} \left(\frac{g_{a}^{4}}{4} \right) \left(\ln(g_{a}^{2}) + \frac{5}{3} \right) \right] \\ + \ln[(t + \bar{t})|\eta(t)|^{4} \left[\left(\sum_{a} \gamma_{i}^{a} \left(\frac{g_{a}^{4}}{4} \right) + \frac{1}{3} \sum_{jk} \gamma_{i}^{jk} k_{s} k_{\bar{s}} \right) \right] \right\},$$

$$(41)$$

 $A_{ijk} = m_{3/2} \bigg\{ (\gamma_i + \gamma_j + \gamma_k) [1 + 2t \cos \theta G_2] \bigg\}$

$$+\frac{\sqrt{3}\sin\theta}{k_{s\bar{s}}^{1/2}} \bigg[k_{s} + \sum_{a} \frac{g_{a}^{2}}{2} (\gamma_{i}^{a} + \gamma_{j}^{a} + \gamma_{k}^{a}) (1 - \ln(g_{a}^{2})) \\ -\ln[(t+\bar{t})|\eta(t)|^{4}] \bigg(\sum_{a} g_{a}^{2} (\gamma_{i}^{a} + \gamma_{j}^{a} + \gamma_{k}^{a}) - \sum_{lm} k_{s} (\gamma_{i}^{lm} + \gamma_{j}^{lm} + \gamma_{k}^{lm}) \bigg) \bigg] \bigg\}.$$

$$(42)$$

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: $M_{\tilde{E},ii}^0$, $M_{\tilde{D},ii}^0$, $M_{\tilde{U},ii}^0$, $M_{\tilde{U},$

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 [34]. 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., $\Gamma(\tilde{\chi}_1^+ \rightarrow \tilde{\chi}_1^0 v l^+) \times BR(W^+ \rightarrow v l^+)$ can be quite different from $\Gamma(\tilde{\chi}_1^+ \rightarrow \tilde{\chi}_1^0 v l^+)$ if the decay $\tilde{\chi}_1^+ \rightarrow \tilde{\chi}_1^0 W^+$ has only small phase space.

The following sfermion decays are calculated:

$$\tilde{f}_i \to f \,\tilde{\chi}_k^0, \, f' \tilde{\chi}_l^\pm, \tag{43}$$

$$\tilde{f}_i \to \tilde{f}_j Z^0, \ \tilde{f}'_j W^{\pm},$$
(44)

$$\tilde{f}_i \to \tilde{f}_j(h^0, H^0, A^0), \ \tilde{f}'_j W^{\pm}.$$
 (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 [35–37]:

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

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

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

where $l = e, \mu, \tau$. The corresponding widths are calculated within SPheno using the formulas given in [36]. In case of GMSB models scenarios exist where the charged sleptons are next to lightest supersymmetric particles (NLSP) and the gravitino \tilde{G} is the LSP. In this case the sleptons decay according to:

$$\tilde{l}_i \to l\tilde{G}.$$
(49)

Here we use the formulas given in [27].

It is well known that the partial widths of sfermions can receive considerable radiative corrections [38]. However, the branching ratios are not that strongly affected [39]. 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:

$$\tilde{\chi}_i^0 \to Z^0 \tilde{\chi}_j^0, \ W^{\pm} \tilde{\chi}_k^{\mp}, \tag{50}$$

$$\chi_i^{\circ} \to (h^{\circ}, H^{\circ}, A^{\circ})\chi_j^{\circ}, \ H^{\perp}\chi_k^{\perp}, \tag{51}$$
$$\tilde{\chi}_i^{0} \to f \,\tilde{f}, \quad \bar{f} \,\tilde{f}. \tag{52}$$

$$\begin{aligned} \chi_i &\to J J_j, \ J J_j, \\ \tilde{\chi}_k^+ &\to Z^0 \tilde{\chi}_s^+, \ W^+ \tilde{\chi}_j^0, \end{aligned} \tag{53}$$

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

$$\tilde{\chi}_k^+ \to f \tilde{f}_i'. \tag{55}$$

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

$$\tilde{\chi}_i^0 \to f \bar{f} \tilde{\chi}_j^0, \ f f' \tilde{\chi}_k^{\mp}, \tag{56}$$

$$\tilde{\chi}_i^0 \to q\bar{q}\tilde{g}, \tag{57}$$

$$\tilde{\chi}_k^+ \to f f \tilde{\chi}_s^+, \ f f' \tilde{\chi}_j^0, \tag{58}$$

$$\tilde{\chi}_k^+ \to q q' \tilde{g}. \tag{59}$$

In the calculation we have included all contributions from gauge bosons, sfermions and Higgs bosons [34,40]. The Higgs contributions can be important in certain regions of parameter space [41]. 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

$$\tilde{\chi}_1^0 \to \gamma \widetilde{G},\tag{60}$$

$$\tilde{\chi}_1^0 \to Z^0 \tilde{G}, \tag{61}$$

$$\tilde{\chi}_1^0 \to h^0 \tilde{G}. \tag{62}$$

Here we use the formulas given in [27].

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

$$\tilde{g} \to q\tilde{q}_i$$
 (63)

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

$$\tilde{g} \to \tilde{\chi}_i^0 q \bar{q},$$
(64)

$$\tilde{g} \to \tilde{\chi}_j^{\pm} q' \bar{q},$$
(65)

$$\tilde{g} \to \bar{b}W^- \tilde{t}_1, \ bW^+ \tilde{t}_1^*.$$
(66)

Here we have implemented the formulas given in [42].

In case of Higgs bosons the following decays are calculated:

 $\phi \to f\bar{f},\tag{67}$

$$\phi \to f_i f_j, \tag{68}$$

$$\phi \to \tilde{\chi}_k^0 \tilde{\chi}_l^0, \tag{69}$$

$$\phi \to \tilde{\chi}_r^+ \tilde{\chi}_s^-,\tag{70}$$

$$H^0 \to Z^0 Z^0, \ W^+ W^-,$$
 (71)

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$$H^0 \to h^0 h^0, \tag{72}$$

$$A^0 \to h^0 Z^0, \tag{73}$$

$$H^+ \to f \bar{f}',$$
 (74)

$$H^+ \to \tilde{f}_i \tilde{f}'_i, \tag{75}$$

$$H^+ \to \tilde{\chi}_k^0 \tilde{\chi}_s^-, \tag{76}$$

$$H^+ \to h^0 W^+ \tag{77}$$

with $\phi = h^0$, H^0 , A^0 and $f = v_i$, $e, \mu, \tau, u, d, c, s, t, b$. It is well known, that the widths as well as the branching ratios of the Higgs bosons can receive large one-loop corrections [43–45]. In the present version only the gluonic QCD corrections for the decays into quarks [43] have been implemented. Therefore, the numbers provided by SPheno have to be taken with care and for refined analysis other programs, such as HDECAY [46] should be used.

4. Production of supersymmetric particles and Higgs bosons

The program calculates the following cross sections:

$$e^+e^- \to f_i f_j \quad (f = l, \nu, q),$$
(78)

$$e^+e^- \to \tilde{\chi}^0_k \tilde{\chi}^0_n, \tag{79}$$

$$e^+e^- \to \chi_r^+ \chi_s^-, \tag{80}$$

$$e^+e^- \to h^0 Z, \ H^0 Z, \tag{81}$$

$$e^+e^- \to h^0 A^0 - H^0 A^0 \tag{82}$$

$$e^+e^- \rightarrow H^+H^-.$$
(82)
(83)

We haven taken the formulas of [47] for sfermion production, [34,48] for production of charginos and neutralinos and [45] for Higgs boson production. Initial state radiation has been included using the formula given in [49]. In case of squarks in addition QCD corrections due to gluon exchange are included [49,50]. 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 [51] 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 $b \rightarrow s\gamma$, the anomalous magnetic moment of the muon a_{μ} and the supersymmetric contributions to the ρ parameter. These constraints are implemented in the program using the formulas given in [52,53] for $b \rightarrow s\gamma$ supplemented by the QCD corrections as given in [54,55] for a_{μ} and [56] for the sfermion contributions to the ρ 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 [57,58] for the case of $b \rightarrow s\gamma$. 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 and gauge boson pole masses, the fine structure constant α , the Fermi constant G_F and the strong coupling constant $\alpha_s(m_Z)$. It is assumed that $\alpha(m_Z)$ and $\alpha_s(m_Z)$ are given in the \overline{MS} scheme.

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. Afterwards 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 [24]. 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 [24]. In case of third generation Yukawa couplings we use the following improvements. In case of the bottom quark we use 3-loop relationship between the pole mass and the \overline{MS} mass $m_{b,\overline{MS}}(m_b)$ as given in [59]. The three-loop RGEs [60] are solved numerically to obtain $m_{b,\overline{MS}}(m_Z)$. Afterwards the shift from the \overline{MS} scheme to the \overline{DR} scheme arising from massless gauge bosons is calculated using the formula [61,62]:

$$m_{b',\overline{DR}}(m_Z) = \left(1 - \frac{\alpha_s}{3\pi} - \frac{35\alpha_s^2}{72\pi^2}\right) m_{b,\overline{MS}}(m_Z),\tag{84}$$

where α_s is given in the \overline{DR} scheme which is the reason for the different factor in front of α_s^2 compared to [62]. We use the complete formulas given in Appendix D of [24] 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 resummed using [63]

$$m_{b,\overline{DR}}(m_Z) = \frac{m_{b',\overline{DR}}(m_Z)}{1 - \frac{\Delta m_{b,\text{SUSY}}(m_Z)}{m_{b,\overline{DR}}(m_Z)}}.$$
(85)

The analogue resumation is done in case of the tau Yukawa coupling. In case of the top quark we take the complete formulas given in Appendix D of [24]. We also add the 2-loop contribution due to gluons in the \overline{DR} scheme [61] which reads as:

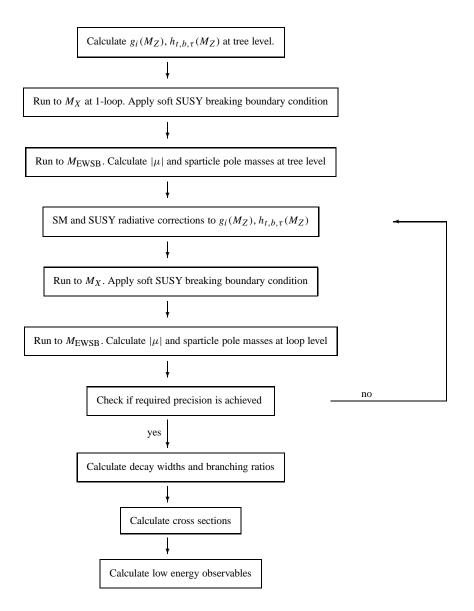


Fig. 1. Algorithm used to calculate the SUSY spectrum, decay widths, production cross sections and low energy observables. Each step (represented by a box) is explained in the text. The initial step is the uppermost one. $M_{\rm EWSB}$ is the scale at which the EWSB conditions are imposed, and M_X is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

$$\Delta m_t^{2\text{-loop}}(Q) = -m_t \alpha_s^2 \left(0.538 - \frac{43}{24\pi^2} L + \frac{3}{8\pi^2} L^2 \right),\tag{86}$$

where m_t is the on-shell top mass and $L = \ln(m_t^2/Q^2)$. The obtained gauge and Yukawa couplings are evolved to the high scale using two-loop RGEs [23]. 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 Section 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 Section 2.3; see also B.3. The complete set of two-loop RGEs [23] is used to evolve the parameters down to the electroweak scale $M_{\text{EWSB}} = \sqrt{m_{\tilde{t}_1} m_{\tilde{t}_2}}$ (there exists also the possibility to set M_{EWSB} by hand as explained 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 [24] for the one-loop contributions and for the 2-loop corrections $O(\alpha_s \alpha_t + \alpha_t^2 + \alpha_s \alpha_b)$ for the neutral Higgs boson the formulas given in [25,26]. 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 $M_{\rm EWSB}$. Also tan β and the vacuum expectation values are evaluated at $M_{\rm EWSB}$ 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 [24]. In all cases running masses are used as input for the loop integrals. In addition we have implemented the $O(\alpha_s \alpha_t + \alpha_t^2 + \alpha_s \alpha_b)$ corrections for the calculation of $|\mu|$ [26]. The numerical evolution of the one-loop integrals is based on the FF package [64] and the LoopTools package [65].

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 δ :

$$\delta > \frac{|m_i - m_{i-1}|}{m_i} \tag{87}$$

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.

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 $M_{\rm EWSB}$. 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.

Afterwards 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 centre 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: $BR(b \rightarrow s\gamma)$, SUSY contributions to a_{μ} and the sfermion contributions to $\Delta \rho$. For 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 $BR(b \rightarrow s\gamma)$ the most important contributions to the C_7 coefficients are implemented as

$$C_7(W^+) = -\frac{K_{ts}K_{tb}x_tW}{4m_W^2} \left(\frac{2}{3}F_1(x_tW) + F_2(x_tW)\right),\tag{88}$$

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$$C_{7}(H^{+}) = -\frac{K_{ts}K_{tb}}{4m_{H^{+}}^{2}} \left[\frac{Y_{t}^{2}\cos^{2}\beta}{4} \left(\frac{2}{3}F_{1}(x_{tH^{+}}) + F_{2}(x_{tH^{+}}) \right) - \frac{Y_{b}Y_{t}\cos\beta\sin\beta m_{t}}{m_{b}} \left(\frac{2}{3}F_{3}(x_{tH^{+}}) + F_{4}(x_{tH^{+}}) \right) \right],$$
(89)

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

$$(90)$$

$$C_{L,ij} = Y_b R_{\tilde{t},i1} U_{j2},\tag{91}$$

$$C_{R,ij} = -gR_{\tilde{t},i1}V_{j1} + Y_tR_{\tilde{t},i2}V_{j2}.$$
(92)

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 [52]. 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 [54] to obtain

$$BR(b \to s\gamma) = 1.258 + 0.382r7^2 + 0.015r_8^2 + 1.395r_7 + 0.161r_8 + 0.083r_7r_8,$$
(93)

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 β effects in case of $b \rightarrow s\gamma$ [58].

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. Afterwards a series of initialization routines are called:

```
Call InitializeControl(11,"SPheno.out","SPheno")
Call InitializeLoopFunctions
Call InitializeStandardModel
```

The routine InitialzeControl opens the file Messages.out at channel 10 where all warnings and/or debugging informations are stored. In addition it reads the file Control.in if present. For the description of Control.in see B.1. Moreover, it opens the output file SPheno.out at channel 11. The last entry "SPheno" is the name of the calling program and is used for the initialization of the debugging system. The routine InitializeLoopFunctions initializes the routines for the calculation of the loop functions which are needed to calculated the loop corrected masses. The routine InitializeStandardModel reads data from the file StandardModel.in. In case that this file is not present, default values are used as described in B.4.

Afterwards the routine HighScaleInput is called:

```
Call HighScaleInput
```

The source code for this routine is given in the file SPheno.f90. This routine reads in the data from the file HighScale.in which specifies the model used as well as the high scale boundary conditions. A short description of the implemented models is given in Section 2.3 and details of the file HighScale.in are given in B.3. Afterwards a first rough estimate of the parameters and couplings is given by calling the routine FirstGuess:

Table 1
Variables for parameters and couplings. The parameters are explained in Section 2. dp means double precision

Parameter/coupling	Type and 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	<pre>complex(dp), dimension(3,3) :: M2_E, M2_L</pre>
M_D^2, M_Q^2, M_U^2	<pre>complex(dp), dimension(3,3) :: M2_D, M2_Q, M2_U</pre>
A_l, A_d, \tilde{A}_u	<pre>complex(dp), dimension(3,3) :: A_1, A_d, A_u</pre>
μ	complex(dp) :: 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	<pre>complex(dp), dimension(3,3) :: Y_1, Y_d, Y_u</pre>
v_1, v_2	real(dp) :: vevSM(2)
g', g, g_s	real(dp) :: gauge(3)

Call FirstGuess(phase_mu, tanb, Mi, M2_E, M2_L, A_l, M2_D & & , M2_Q, M2_U, A_d, A_u, mu, B, M2_H, gp, g, Y_l & & & , Y_d, Y_u, vevSM, mP02, mP0)

The meaning of the various variables and their type is given in Table 1 except for mP02, mP0 which are given in Table 2. The parameters obtained are used to get the initial values for the spectrum at tree-level:

The variables for the masses and mixing matrices are given in Table 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 6×6 (except for sneutrions which is a 3×3 matrix). In the present release most of the entries are zero except for the diagonal 2×2 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.

These initial values serve as input for the more accurate calculation of the parameters including higher oder corrections. Before calling the subroutine Sugra, which performs these calculations, 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 $M_{\text{EWSB}} = \sqrt{m_{\tilde{t}_1} m_{\tilde{t}_2}}$.

Masses/mixing matrix	Type and Fortran name
m _ğ	real(dp) :: mglu
$e^{\tilde{\varphi_{g}}}$	complex(dp) :: PhaseGlu
$m_{\tilde{\chi}_i^+}$	real(dp) :: mC(2)
U, V	complex(dp) :: U(2,2), V(2,2)
$m_{\tilde{\chi}_{j}^{0}}$	real(dp) :: N(4)
N	complex(dp) :: N(4,4)
m_{h^0}, m_{H^0}	real(dp) :: mS0(2)
R_{α}	real(dp) :: RS0(2,2)
m_{G^0}, m_{A^0}	real(dp) :: mP0(2)
R_{β}	real(dp) :: RP0(2,2)
$m_{G^{+}}, m_{H^{+}}$	real(dp) :: mSpm(2)
R'_{β}	complex(dp) :: RSpm(2,2)
$m_{\tilde{v}}$	real(dp) :: mSneut(3)
$R_{ ilde{ u}}$	complex(dp) :: Rsneut(3,3)
$m_{\tilde{l}}$	real(dp) :: mSlepton(6)
$R_{\tilde{l}}$	complex(dp) :: Rslepton(6,6)
$m_{\tilde{u}}$	real(dp) :: mSup(6)
$R_{\tilde{u}}$	complex(dp) :: Rsup(6,6)
$m_{\tilde{d}}$	real(dp) :: mSdown(6)
$R_{\tilde{d}}^{u}$	complex(dp) :: Rsdown(6,6)
U_u^L, U_u^R	complex(dp) :: uU_L(3,3), uU_L(3,3)
U_d^L, U_d^R	complex(dp) :: uD_L(3,3), uD_L(3,3)
$U_l^{\tilde{L}}, U_l^{\tilde{R}}$	complex(dp) :: uL_L(3,3), uL_L(3,3)

Variables for masses and mixing matrices as given by the routine Sugra.
Their connection to the parameters at tree-level is explained in Section 2. dp
means double precision

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

```
delta = 1.e-3_dp
WriteOut = .False.
n_run = 20
Call Sugra(delta, 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, uL_L, uL_R, uD_L, uD_R, uU_L, uU_R, 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, m_GUT, kont, WriteOut, n_run)
```

The meaning of the various variables and their type is given in Tables 1 and 2. Note that the quantities uL_L , uL_R , uD_L , uD_R , uU_L , uU_R are 3×3 unit matrices in the present implementation. Also all SUSY parameters, Yukawa couplings and the sfermion masses are flavor diagonal. However, we have constructed the interface already in such a way that an extension to include mixing effects between the generations is facilitated. Compared to the previous calls of routines the following new variable 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 color 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 1loop level using the formulas given in [24]. The exceptions are the masses of the neutral Higgs bosons (scalar and pseudo-scalar), where in addition the 2-loop corrections $O(\alpha_s \alpha_t + \alpha_t^2 + \alpha_s \alpha_b)$ are included [25,26], and the μ parameter where also the $O(\alpha_s \alpha_t + \alpha_t^2 + \alpha_s \alpha_b)$ corrections [26] are included.

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 = 1.e-4_dp
 If (HighScaleModel.Eq."GMSB") Is_GMSB = .True.
 Call CalculateBR(gauge, mGlu, PhaseGlu, mC, U, V, mN, N
                                                            8
    & , mSneut, RSneut, mSlepton, RSlepton, mSup, RSup
                                                            &
    & , mSdown, RSdown, uL_L, uL_R, uD_L, uD_R, uU_L, uU_R &
    & , mS0, RS0, mP0, RP0, mSpm, RSpm, epsI, deltaM
                                                            &
    & , CalcTBD, kont, ratioWoM, Y_d, A_d, Y_l, A_l, Y_u
                                                            &
    & , A_u, mu, vevSM, Fgmsb, m32
                                                            8
    & , 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_C2, gT_C
                                                            &
    & , BR_C2, gP_C3, BR_C3, gP_N2, gT_N, BR_N2, gP_N3
                                                            &
    & , BR_N3, gP_G2, gT_G, BR_G2, gP_G3, BR_G3, gP_P0
                                                            &
    & , gT_P0, BR_P0, gP_S0, gT_S0, BR_S0, 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:

- 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.

1		1	ξ, β	
Mode	ĩ	ũ	ã	ũ
$\tilde{f}_i \to f \tilde{\chi}_k^0$	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
$\tilde{f}_i \to f \tilde{g}$	-	-	7	7
$\tilde{f}_i \to W^{\pm} \tilde{f}'_i$	7	7–8	8–9	8–9
$\tilde{f}_i \to H^{\pm} \tilde{f}'_i$	8	9–10	10-11	10-11
$\tilde{f}_2 \rightarrow Z^0 \tilde{f}_1$	9	11	12	12
$\tilde{f}_2 \to A^0 \tilde{f}_1$	10	12	13	13
$\tilde{f}_2 \rightarrow h^0 \tilde{f}_1$	11	13	14	14
$\tilde{f}_2 \to H^0 \tilde{f}_1$	12	14	15	15
$\tilde{l}_1 \rightarrow l \widetilde{G}$	13	_	_	-

Table 3			
Correspondence between the indices for sfermion p	partial widths (bra	anching ratios)	and the modes

Correspondence between the second indices for chargino partial widths (branching ratios) and the decay modes. In case of two body decays the variables are gP_C2 (BR_C2) and in case of three body decays gP_C3 (BR_C3)

Mode	Index of gP_C2 (BR_C2)	Index of gP_C3 (BR_C3)
$\tilde{\chi}_i^+ \rightarrow \tilde{l}_{m,k}^+ \nu_m$	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	
$\tilde{\chi}^+_i ightarrow \tilde{\chi}^0_j W^+$	22–25	
$\tilde{\chi}_i^+ \rightarrow \tilde{\chi}_i^0 H^+$	26–29	
$\tilde{\chi}_2^+ \rightarrow \tilde{\chi}_1^+ Z^0$	30	
$\tilde{\chi}_2^+ \rightarrow \tilde{\chi}_1^+ A^0$	31	
$\tilde{\chi}_2^+ \rightarrow \tilde{\chi}_1^+ h^0$	32	
$\tilde{\chi}_2^+ \rightarrow \tilde{\chi}_1^+ H^0$	33	
$\tilde{\chi}_i^+ \to \tilde{\chi}_j^0 u_m \bar{d}_m$		1–12
$\tilde{\chi}_i^+ \to \tilde{\chi}_i^0 l_m^+ \nu_m$		13–24
$\tilde{\chi}_i^+ \to \tilde{g} u_m \bar{d}_m$		25–27
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ u_m \bar{u}_m$		28-30
$\tilde{\chi}_2^+ \to \tilde{\chi}_1^+ d_m \bar{d}_m$		31–33
$\tilde{\chi}_2^+ \rightarrow \tilde{\chi}_1^+ l_m l_m^+$		34–36
$\tilde{\chi}_2^+ \rightarrow \tilde{\chi}_1^+ \sum_m v_m \bar{v}_m$		37

- 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.
- 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

Table 5

Correspondence between the second indices for neutralino partial widths (branching ratios) and the decay modes. In case of two body decays the variables are gP_C2 (BR_C2) and in case of three body decays gP_C3 (BR_C3)

Mode	Index of gP_C2 (BR_C2)	Index of gP_C3 (BR_C3)
$\tilde{\chi}_i^0 \rightarrow \tilde{l}_{m\ k}^+ l_m$	1–6	
$\tilde{\chi}_i^0 \to \tilde{\nu}_m l_m^+$	7–9	
$\tilde{\chi}_i^0 \to \bar{\tilde{u}}_{m,k} u_m$	10–15	
$\tilde{\chi}_i^0 \to \tilde{d}_{m,k} \bar{d}_m$	16–21	
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^- W^+$	22–23	
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^- H^+$	24–25	
$\tilde{\chi}_i^0 \to \tilde{\chi}_i^0 Z^0$	26-(24+i)	
$\tilde{\chi}_i^0 \to \tilde{\chi}_i^0 A^0$	(25+i)-(23+2i)	
$\tilde{\chi}_i^0 \rightarrow \tilde{\chi}_i^0 h^0$	(26+i)-(22+3i)	
$\tilde{\chi}_i^0 \to \tilde{\chi}_i^0 H^0$	(27+i)-(21+4i)	
$\tilde{\chi}_1^0 \to \gamma \widetilde{G}$	74	
$\tilde{\chi}_1^0 \to Z^0 \tilde{G}$	75	
$\tilde{\chi}_1^0 \to h^0 \widetilde{G}$	76	
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^0 u_m \bar{u}_m$		1–(3*i – 3)
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^0 d_m \bar{d}_m$		(3*i - 2)–(6*i - 6)
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^0 l_m^+ l_m^-$		(6*i - 5)–(9*i - 9)
$\tilde{\chi}_i^0 \to \tilde{\chi}_j^0 \sum_n \nu_m \bar{\nu}_m$		(9*i - 8)-(10*i - 9)
$\tilde{\chi}_i^0 \to \widetilde{G} u_m \bar{u}_m$		(10*i - 8)–(10*i - 6)
$\tilde{\chi}_i^0 \to \widetilde{G} d_m \bar{d}_m$		(10*i - 5)–(10*i - 3)

Correspondence between the indices for gluino partial widths (branching ratios) and the decay modes. In case of two body decays the variables are gP_G2 (BR_G2) and in case of three body decays gP_G3 (BR_G3)

Mode	Index of gP_C2 (BR_C2) Index of gP_C3 (BR_C2)	
$\tilde{g} \to \bar{\tilde{d}}_{m,k} d_m$	1–6	
$\tilde{g} \to \tilde{u}_{m,k} \bar{u}_m$	7–12	
$\tilde{g} \to \tilde{\chi}_{j}^{0} u_{m} \bar{u}_{m}$		1–12
$\tilde{g} \rightarrow \tilde{\chi}_{i}^{0} d_{m} \bar{d}_{m}$		13–24
$\tilde{g} \to \tilde{\chi}_i^+ d_m \bar{u}_m$		25-30
$\tilde{g} \rightarrow \tilde{t}_i W^- \bar{b}$		31–32
$\tilde{g} \rightarrow \tilde{t}_1 \bar{c}$		33

ratio of the width Γ over the mass *m* of the decaying particle is small: $\Gamma/m < \texttt{ratioWoM}$, then three body decay modes are calculated.

• 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.

Table /	Tabl	e	7
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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_S0 (BR_S0) 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 ϕ stands for h^0 , H^0 and A^0 . The index runs from 1 to 3

Mode	h^0	H^0	A^0
$\phi \to l_m^+ l_m^-$	1–3	1–3	1–3
$\phi \to d_m \bar{d}_m$	4–6	4–6	4–6
$\phi \rightarrow u_m \bar{u}_m$	7–9	7–9	7–9
$H^0 \rightarrow \tilde{l}^+_{m,1} \tilde{l}^{m,1}$	-	$10 + 3^{*}(m - 1)$	_
$\phi \to \tilde{l}_{m,1}^{\mp} \tilde{l}_{m,2}^{\pm}$	_	$11 + 3^*(m-1)$	$11 + 3^*(m-1)$
$H^0 \rightarrow \tilde{l}^+_{m,2} \tilde{l}^{m,2}$	_	$12 + 3^*(m-1)$	_
$H^0 \to \tilde{\nu}_m \overline{\tilde{\nu}}_m$	-	19 + m	_
$H^0 \rightarrow \tilde{d}_{m,1} \overline{\tilde{d}}_{m,1}$	_	22 + 3*(m - 1)	_
$\phi \to \tilde{d}_{m,1}\overline{\tilde{d}}_{m,2}$	_	23 + 3*(m - 1)	20 + 3*(m-1)
$H^0 \rightarrow \tilde{d}_{m,2} \overline{\tilde{d}}_{m,2}$	-	24 + 3*(m - 1)	_
$H^0 \to \tilde{u}_{m,1} \overline{\tilde{u}}_{m,1}$	-	31 + 3*(m - 1)	_
$\phi \to \tilde{u}_{m,1} \overline{\tilde{u}}_{m,2}$	-	32 + 3*(m - 1)	29 + 3*(m-1)
$H^0 \to \tilde{u}_{m,2}\overline{\tilde{u}}_{m,2}$	_	33 + 3*(m - 1)	_
$\phi \to \tilde{\chi}^0_r \tilde{\chi}^0_s \; (r \leqslant s)$	40–49	40–49	37–46
$\phi \to \tilde{\chi}_k^+ \tilde{\chi}_n^- \ (k \leq n)$	50-52	50-52	47–49
$H^0 \rightarrow Z^0 Z^0$	_	53	_
$H^0 \rightarrow W^+ W^-$	-	54	-
$H^0 \rightarrow h^0 h^0$	-	59	-
$A^0 \rightarrow h^0 Z^0$	-	_	52

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

Process	Fortran name and type
$e^+e^- \rightarrow \tilde{u}_i \tilde{u}_j$	real(dp) :: SigSup(6,6)
$e^+e^- \rightarrow \tilde{d}_i \tilde{d}_j$	real(dp) :: SigSdown(6,6)
$e^+e^- \rightarrow \tilde{l}_i \tilde{l}_j$	real(dp) :: SigSle(6,6)
$e^+e^- \rightarrow \tilde{v}_i \tilde{v}_j$	real(dp) :: SigSn(6,6)
$e^+e^- \rightarrow \tilde{\chi}^0_k \tilde{\chi}^0_n$	real(dp) :: SigChi0(4,4)
$e^+e^- \rightarrow \tilde{\chi}_r^+ \tilde{\chi}_s^-$	real(dp) :: SigN(4,4)
$e^+e^- \rightarrow h^0 Z, \ H^0 Z$	real(dp) :: SigS0(2)
$e^+e^- \rightarrow h^0 A^0, \ H^0 A^0$	real(dp) :: SigSP(2)
$e^+e^- \rightarrow H^+H^-$	real(dp) :: SigHp

Correspondence between the production cross sections and the variables used in the program

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

```
If ((L_CS).and.(kont.eq.0)) then
 Call InitializeCrossSections(Ecms, Pm, Pp, ISR)
 Call CalculateCrossSections(Ecms, Pm, Pp, ISR
                                                               &
          & , mSup, RSup, mf_u, mSdown, RSdown, mf_d, mglu
                                                               8
          & , SigSup, SigSdown, mSlepton, RSlepton
                                                               &
          & , mSneut, RSneut, SigSle, SigSn, mC, U, V, mN, N
                                                               &
          & , SigC, SigChiO, mSO, RSO, vevSM, mPO, RPO, mSpm
                                                               8
          & , RSpm, SigSO, SigSP, SigHp )
End If
```

Table 9

Here the following additional input is needed:

- Ecms: the centre 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 [49].

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 2×2 diagonal blocks and they are sorted according to the generations as in the case of the sfermion mixing matrices.

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

```
If (kont.eq.0) then
Call CalculateLowEnergyConstraints(gauge, Y_l, Y_d, Y_u
                                                             &
  & , mSpm2, RSpm, mC, U, V, mN, N, mSup2, RSup, mSdown2
                                                             &
  & , RSdown, mSlepton2, RSlepton, mSneut2, RSneut
                                                             &
    , BRbtosgamma, a_mu, Delta_Rho)
   &
Else
```

```
BRbtosgamma = 0._dp
a_mu = 0._dp
Delta_Rho = 0._dp
End If
```

Here BRbtosgamma, a_mu, and Delta_Rho denote $10^4 \times BR(b \rightarrow s\gamma)$, the SUSY contributions to a_{μ} and the sfermion contributions to $\Delta \rho$, respectively.

Afterwards the statement

```
Call WriteOutPut0(11, 1.e-6_dp, 1.e-3_dp)
```

is used to write all information to the file connected with unit 11 (first entry). The second entry puts a lower bound on the branching ratios to be written. In the case above, branching ratios smaller than 10^{-6} will not be given. The third entry gives the minimum value for the cross section in fb which will be written to the output file. In the example given above cross sections smaller than 10^{-3} will not be written to the output file.

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 μ 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: $BR(b \rightarrow s\gamma)$, the supersymmetric contributions to the anomalous magnetic moment of the muon a_{μ} and the sfermion contributions to the parameter.

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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Appendix A. Switches

In this appendix we describe the switches for influencing the behavior 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 $M_{\text{EWSB}} = \sqrt{m_{\tilde{l}_1} m_{\tilde{l}_2}}$. By calling

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

 $M_{\rm EWSB}$ 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 [66]. 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..

Appendix 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.

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 file Control.in is not present. Here ErrorLevel is an integer in the range [-2, 2] where the numbers correspond to the following behavior 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 electrons
0. ! 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 file CrossSections.in is not present.

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 [67]. In Appendix C 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 μ 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 β . The file reads, for example,

msugra		
250.	!	M_1/2
100.	!	M_0
-100.	!	A_0
10.	!	tan(beta)
1.	!	sign of mu
.TRUE.	!	if 2-loop RGEs should be used

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_{ν_R} and the light neutrino masses m_{ν_i} (*i* = 1, 2, 3). The file reads, for example,

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

B.3.3. GMSB

GMSB

The GMSB scenario is characterized by two mass parameters M_M and Λ ; 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 μ as described in Section 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,

CLOID		
100000.	! Lambda	
200000.	! M_M	
1	! N_5	
0	! N_10	
0.	! A0	
15.	! tan(beta)	
1.	! sign of mu	
.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 β and the sign of μ . 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 $\langle t \rangle$ 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 δ_{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 β and the sign of μ . The file reads, for example, as

String_	OI
---------	----

180.	!	!	M_3/2
14.0	!	!	<t></t>
0.5	!	!	g_s^2
0.9	!	!	<pre>sin^2(theta)</pre>
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 $\langle t \rangle$, 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 counter-term δ_{GS} . Moreover, one needs to specify $\tan \beta$ and the sign of μ . There are two different scenarios implemented denoted as String_OIIa and String_OIIb corresponding to boundary conditions (*A*) and (*B*) of [31], respectively. The files read, for example, as

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

and

String_011b	
300. !	M_3/2
14.6 !	<t></t>
0.5 !	g_s^2
0.9 !	<pre>sin^2(theta)</pre>
0. !	delta_GS
5. !	tan(beta)
1. !	phase(mu)
.TRUE. !	if 2-loop RGEs should be used

B.3.7. SUGRA

Church of Th

This input files 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: $M^0_{\tilde{E},ii}$, $M^0_{\tilde{D},ii}$, $M^0_{\tilde{U},ii}$, $M^0_{\tilde{Q},ii}$; two Higgs mass parameters $M^0_{H_1}$ and $M^0_{H_2}$; 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	a			
480.	300.	300.	!	M_1/2_i
150.	150.	150.	!	MO_E_ii
150.	150.	150.	!	MO_L_ii
150.	150.	150.	!	M0_D_ii
150.	150.	150.	!	M0_Q_ii

150. 150. 150.	!	MO_U_ii
150. 150.	!	MO_H_i
0. 0. 0.	!	A0_u_ii
0. 0. 0.	!	A0_d_ii
0. 0. 0.	!	A0_e_ii
10.	!	tan(beta)
1.	!	phase(mu)
.TRUE.	!	if 2-loop RGEs should be used

B.4. StandardModel.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:

```
91.187
                          ! Z boson mass
2.490
                         ! width of Z-boson
0.14 0.14 0.14 0.14 0.14 ! Br(Z -> q q) q=u,d,c,s,b
                         ! Br(Z -> 1 1) 1=e, mu, tau
0.035 0.035 0.03
                          ! Br(Z -> invisible)
0.2
80.41
                          ! W boson mass
2.06
                         ! widht of W-boson
0.35 0.35
                         ! Br(W -> q q)
0.1 0.1 0.1
                         ! Br(W -> l nu)
0.51099906e-3
                         ! electron mass
0.105658
                         ! muon mass
1.7771
                         ! tau mass
0.004
                         ! u-quark mass
1.4
                         ! c-quark mass
174.3
                         ! t-quark mass
0.01
                         ! d-quark mass
                         ! s-quark mass
0.15
4.62
                         ! b-quark mass
0.2315
                         ! sin(theta_W) squared
137.0359895
                         ! 1 / Alpha
                         ! 1 / Alpha at m_Z
127.9
0.119
                          ! Alpha_s at m_Z
1.16639e-5
                          ! G_F, Fermi constant
              ! s12 of CKM, parametrized as particle data book
0.220
0.039
              ! s23 of CKM
0.0031
             ! s13 of CKM
Ο.
              ! phase of CKM
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.

Appendix C. 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 B.3.1 and the default values of the files Control.in, CrossSections.in and StandardModel.in.

```
mSugra input at the GUT scale 1.8193751412503776E+16
M_1/2
      :
             2.5000000000000000E+02
             1.0000000000000000E+02
M_0
         :
A_0
        :
             tan(beta) :
           10.00000000000000000
             1.000000000000000000
phase(mu) :
Parameters at the scale
                          4.8387163690920886E+02
Gaugino mass parameters
  1.0256160300628754E+02
                                                  5.8695416971423458E+02
                          1.9230453743751079E+02
tan(beta), mu, B
10.00000000000000000
                          3.5489208425079283E+02 1.5675353981646900E+04
Slepton mass parameters
A 1
 -7.2960260902552873E-03 -1.5085451901504909
                                                -25.2233192083639537
M2 E
                          1.8407069022083171E+04
                                                  1.7758366624396785E+04
 1.8409373246520361E+04
M2_L
  3.7890957725749206E+04
                          3.7889823466403272E+04
                                                  3.7570561651708413E+04
Squark mass parameters
A_d
-0.2294422308452373
                                                 -1.0705947183168641E+02
                         -3.4416136998921720
Αu
 -7.4845500397144014E-03 -2.6195690984405484
                                                 -4.4984080721097047E+02
M2 D
 2.7694850270662527E+05
                        2.7694564371571731E+05
                                                  2.7392874700277165E+05
M2_U
  2.7918011844082910E+05
                          2.7917785219910560E+05
                                                  1.7963017265859758E+05
M2_0
  2.9890503958833648E+05
                          2.9890250621627242E+05
                                                  2.4881821604275642E+05
Higgs mass parameters
  3.2728705451079368E+04 -1.2647888724841128E+05
Masses and mixing matrices
Gluino :
          5.9431308258414890E+02 1.000000000000000
Charginos
  1.8128399897496990E+02
                          3.7972508491362004E+02
 ΤT
  -0.915146173
                0.403122168
  0.403122168
                0.915146173
 v
  -0.969411416
                0.245441450
  0.245441450
                0.969411416
Neutralinos
 97.6301037960382843 1.8286658240938323E+02 3.6169325535757582E+02
 3.7956517886694593E+02
```

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1	3																			
1	4	(0.0																	
2	1	(-0.1						0											
2	2	(-0.9						0						,					
2	3	(0.2						0						,					
2	4	(-0.1						0						,					
3	1	(0.0						0						,					
3	2	(-0.0						-0											
3	3	(-0.0				,		-0						,					
3	4	(-0.0						-0						,					
4	1	(0.1						0_0						- /					
4	2	(
4	3		-0.6												,					
4	4		0.6																	
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tau-s									560											
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```
t-squark masses
  3.9757274604952244E+02 5.8855899352335769E+02
R_t
  0.554536381 0.832159481
 -0.832159481 0.554536381
d-squark masses
 5.4761741237742456E+02 5.6953755385412876E+02
R_d
  0.000836413 0.999999650
  -0.999999650 0.000836413
s-squark masses
 5.4761111206790952E+02 5.6953840728628120E+02
R s
  0.012543035 0.999921333
 -0.999921333 0.012543035
b-squark masses
 5.1664747823588539E+02 5.4745980835840328E+02
Rb
  0.951026325 0.309109898
 -0.309109898 0.951026325
m_A0, m_H+
 3.9716087636273568E+02
                        4.0654002309066368E+02
m_h0, m_H0
 1.1166181009637616E+02 3.9789685848133115E+02
R_S0
  0.113896390 0.993492633
 -0.993492633 0.113896390
Low energy constraints
10^4 Br(b -> s gamma) : 0.3838177E+01
Delta(a_mu)
              : 0.5447498E-08
Delta(rho)
                    : 0.1829291E-03
                  Decay widths (GeV) and branching ratios
Selectron_1
 Neutralino_1 e
                             0.21131224 100.0000000
Total width :
                             0.21131224
Selectron_2
 Neutralino_1 e
                                         52.26317145
                            0.12712238
 Neutralino_2 e
                             0.04013733 16.50145248
 Chargino_1 neutrino
                         0.07597540 31.23537607
 Total width :
                            0.24323511
```

Smuon_1		
Neutralino_1 mu	0.21104744	100.00000000
Total width :	0.21104744	
Gmuon 0		
Smuon_2	0.12728238	52.28029028
Neutralino_1 mu Neutralino_2 mu		16.49625242
	0.04016202 0.07601710	31.22345730
Chargino_1 neutrino Total width :	0.24346150	31.22345/30
iotal width .	0.24340150	
Stau_1		
Neutralino_1 tau	0.14597684	100.00000000
Total width :	0.14597684	100100000000
	012109/001	
Stau_2		
_ Neutralino_1 tau	0.16310412	55.22030445
Neutralino_2 tau	0.04619971	15.64130946
Chargino_1 neutrino	0.08606600	29.13838609
Total width :	0.29536983	
e-Sneutrino		
Neutralino_1 neutrino	0.15829938	90.39621196
Neutralino_2 neutrino	0.00372885	2.12934196
Chargino_1 e	0.01308905	7.47444608
Total width :	0.17511727	
mu-Sneutrino		
Neutralino_1 neutrino	0.15829302	90.40278693
Neutralino_2 neutrino	0.00372587	2.12788095
Chargino_1 mu	0.01307861	7.46933212
Total width :	0.17509749	
tau-Sneutrino		
Neutralino_1 neutrino	0.15649412	92.23051574
Neutralino_2 neutrino	0.00292840	1.72586668
Chargino_1 tau	0.01025464	6.04361757
Total width :	0.16967717	
Sdown_1		
Neutralino_1 d-quark	0.28784101	98.48847777
Neutralino_2 d-quark	0.00280074	0.95830889
Neutralino_2 d-quark Neutralino_3 d-quark	0.00038028	0.13011842
Neutralino_4 d-quark	0.00123347	0.42204919
Chargino_1 u-quark	0.0000306	0.00104559
Total width :	0.29225857	0.00104555
	0.27225057	
Sdown_2		
Neutralino_1 d-quark	0.12768972	2.43021937
Neutralino_2 d-quark	1.60529969	30.55242392
Neutralino_3 d-quark	0.00884812	0.16839948
Neutralino_4 d-quark	0.08754132	1.66610595
Chargino_1 u-quark	3.18849528	60.68415751

Chargino_2 u-quark Total width :	0.23637246 5.25424660	4.49869376
S-strange_1 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 :	0.28785440 0.00313853 0.00039188 0.00123018 0.00068723 0.00000010 0.29330232	98.14255868 1.07006608 0.13360864 0.41942337 0.23430866 0.00003458
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 :	0.12767855 1.60496370 0.00888876 0.08758857 3.18774099 0.23649323 5.25335380	2.43041977 30.55122040 0.16920168 1.66728859 60.68011250 4.50175707
Sbottom_1 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 :	0.16184398 1.32661019 0.01759582 0.04615410 1.71616072 0.66040512 3.92876992	4.11945668 33.76655332 0.44787087 1.17477226 43.68188400 16.80946286
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 :	0.24397318 0.08601743 0.03700740 0.05304386 0.11582889 0.19366391 0.72953467	33.44230042 11.79072502 5.07274017 7.27091758 15.87709181 26.54622500
Sup_1 Neutralino_1 u-quark Neutralino_2 u-quark Neutralino_3 u-quark Neutralino_4 u-quark Chargino_1 d-quark Total width :	1.15181264 0.01120273 0.00152307 0.00494195 0.00000002 1.16948041	98.48926332 0.95792374 0.13023466 0.42257635 0.00000193
Sup_2 Neutralino_1 u-quark Neutralino_2 u-quark Neutralino_3 u-quark Neutralino_4 u-quark	0.03620866 1.72553982 0.00517093 0.06407609	0.66423008 31.65418423 0.09485826 1.17544453

Chargino_1 d-quark Chargino_2 d-quark Total width :	3.53559056 0.08463583 5.45122190	64.85867992 1.55260297
S-charm_1 Neutralino_1 c-quark Neutralino_2 c-quark Neutralino_3 c-quark Neutralino_4 c-quark Chargino_1 s-quark Chargino_2 s-quark Total width :	1.15098392 0.01253940 0.00158298 0.00492241 0.00276378 0.00000107 1.17279356	98.14036844 1.06919075 0.13497525 0.41971639 0.23565786 0.00009130
S-charm_2 Neutralino_1 c-quark Neutralino_2 c-quark Neutralino_3 c-quark Neutralino_4 c-quark Chargino_1 s-quark Chargino_2 s-quark Total width :	0.03701501 1.72415946 0.00516765 0.06416446 3.53281680 0.08472953 5.44805290	0.67941726 31.64725980 0.09485322 1.17775023 64.84549368 1.55522582
Stop_1 Neutralino_1 t-quark Neutralino_2 t-quark Chargino_1 b-quark Chargino_2 b-quark Total width :	0.39250239 0.22674584 1.38120247 0.02203615 2.02248685	19.40691906 11.21123908 68.29228464 1.08955722
<pre>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 :</pre>	0.22161960 0.62619038 0.34624563 1.54482109 1.54991244 1.51409588 1.51313204 0.27752232 7.59353939	2.91852834 8.24635720 4.55973974 20.34388720 20.41093566 19.93926416 19.92657136 3.65471634
Chargino_1 neutralino_1 d u neutralino_1 s c neutralino_1 e nu neutralino_1 mu nu neutralino_1 tau nu Total width :	0.00033235 0.00033232 0.00058806 0.00063093 0.01474527 0.01662893	1.99860842 1.99844982 3.53637657 3.79415091 88.67241428
Chargino_2 Selectron_2 neutrino Smuon_1 neutrino Smuon_2 neutrino	0.12652636 0.00000150 0.12656625	5.19161956 0.00006154 5.19325658

<pre>Stau_1 neutrino Stau_2 neutrino e-sneutrino e mu-sneutrino mu tau-sneutrino tau Neutralino_1 W Neutralino_2 W Chargino_1 Z Chargino_1 h0 Total width :</pre>	0.00104893 0.13749520 0.05302870 0.05309068 0.07070699 0.17302320 0.67940247 0.58927646 0.42696015 2.43712690	0.04303979 5.64169248 2.17586926 2.17841255 2.90124380 7.09947452 27.87718885 24.17914547 17.51899560
Neutralino_1 : stable		
Neutralino_2 Selectron_1 e Smuon_1 mu Stau_1 tau Total width :	0.00154723 0.00160018 0.01862788 0.02177529	7.34859632
Neutralino_3 Selectron_1 e Selectron_2 e Smuon_1 mu Smuon_2 mu Stau_1 tau Stau_2 tau e-sneutrino neutrino mu-sneutrino neutrino tau-sneutrino neutrino Chargino_1 W Neutralino_1 Z Neutralino_2 Z Neutralino_2 h0 Total width :	0.00503923 0.00218588 0.00508346 0.00227146 0.02025229 0.02419151 0.01234523 0.01234554 0.01243421 1.14827835 0.22657242 0.40462398 0.04071375 0.02112676 1.93746405	
Neutralino_4 Selectron_1 e Selectron_2 e Smuon_1 mu Smuon_2 mu Stau_1 tau Stau_2 tau e-sneutrino neutrino mu-sneutrino neutrino tau-sneutrino neutrino Chargino_1 W Neutralino_1 Z Neutralino_2 Z Neutralino_2 h0 Total width :	0.02090064 0.05133031 0.02086555 0.05145801 0.01565228 0.08419943 0.13486416 0.13486717 0.13571657 1.34016810 0.05567616 0.04779240 0.17892991 0.35146997 2.62389066	0.79655160 1.95626710 0.79521429 1.96113393 0.59652937 3.20895348 5.13985427 5.13996895 5.17234103 51.07560769 2.12189348 1.82143249 6.81925922 13.39499312

Gluino		
	0.32555602	9.88307192
Sup_1 u-quark Sup_2 u-quark	0.13408883	4.07060373
S-charm_1 c-quark	0.32506240	9.86808701
S-charm_2 c-quark	0.13420743	4.07420414
Stop_1 t-quark	0.35279908	10.71010362
Sdown_1 d-quark	0.32809509	9.96015176
Sdown_2 d-quark	0.09594182	2.91255533
S-strange_1 s-quark	0.32815106	9.96185087
S-strange_2 s-quark	0.09594849	2.91275775
Sbottom_1 b-quark	0.82565254	25.06475973
Sbottom_2 b-quark	0.34857445	10.58185414
Total width :	3.29407721	
h0		
electrons	0.0000000	0.00000125
muons	0.00000103	0.05351444
taus	0.00029239	15.14367767
d-quark	0.0000001	0.00031725
s-quark	0.00000138	0.07138010
b-quark	0.00152831	79.15466880
u-quark	0.00000000	0.00004556
c-quark	0.00010767	5.57639494
Total width :	0.00193079	5.5,055151
iotai width ·	0.00193079	
но		
muons	0.00028014	0.04201232
taus	0.07938694	11.90538074
d-quark	0.0000168	0.00025219
s-quark	0.00037836	0.05674169
b-quark	0.42140146	63.19609693
c-quark	0.00000511	0.00076636
t-quark	0.03081125	4.62065503
Selectron 1 1	0.00039974	0.05994791
Smuon 1 1	0.00040732	0.06108426
Smuon 1 2	0.00003734	0.00559954
Stau 1 1	0.00433734	0.65045550
Stau 1 2	0.00799437	1.19888731
e-Sneutrino	0.00073208	0.10978775
mu-Sneutrino	0.00073222	0.10980771
tau-Sneutrino	0.00076881	0.11529554
neutralino_1 neutralino_1	0.01687144	2.53015073
neutralino_1 neutralino_2	0.04686601	7.02833048
neutralino_2 neutralino_2	0.01058042	1.58670780
chargino_1 chargino_1	0.03027976	4.54094909
Z Z	0.00153237	0.22980460
W W	0.00320663	0.48088662
h0 h0	0.00980485	1.47039879
Total width :	0.66681564	1.11059019
iotai wiutii ·	0.00001004	
A0		
muons	0.00028050	0.02631577

taus	0.07949301	7.45791191
d-quark	0.00000168	0.00015796
s-quark	0.00037883	0.03554126
b-quark	0.42198287	39.58978482
c-quark	0.0000389	0.00036525
t-guark	0.10032310	9.41215930
Smuon 1 2	0.00003869	0.00362993
Stau 1 2	0.01130912	1.06100403
neutralino_1 neutralino_1	0.02607822	2.44661842
neutralino_1 neutralino_2	0.10959778	10.28229529
neutralino_2 neutralino_2	0.08682800	8.14606947
chargino_1 chargino_1	0.22688237	21.28575532
h0 Z	0.00269020	0.25239056
Total width :	1.06588826	0.23237030
iotai width ·	1.00500020	
H^+		
electron neutrino	0.0000001	0.0000109
muon neutrino	0.00028712	0.04644904
tau neutrino	0.08137042	13.16371935
d-quark u-quark	0.00000151	0.00024435
s-quark c-quark	0.00034333	0.05554153
b-quark t-quark	0.37287242	60.32152442
Selectron_2 Sneutrino	0.00053557	0.08664129
Smuon_1 Sneutrino	0.00005577	0.00902161
Smuon_2 Sneutrino	0.00053145	0.08597568
Stau_1 Sneutrino	0.01497337	2.42231984
Stau_2 Sneutrino	0.00000770	0.00124494
chargino_1 neutralino_1	0.14336239	23.19248520
chargino_1 neutralino_2	0.00074334	0.12025399
h0 W	0.00305719	0.49457746
Total width :	0.61814157	0.1913//10
10001 112011	0.0101110/	
Total cross sections in fb for	r:	
E_cms: 500.0000000 GeV		
Degree of polarization: P_e-		e + = 0.000000
Intial state radiation is inc		
u-squarks : kinematically not	t possible	
c-squarks : kinematically not	t possible	
t-squarks : kinematically not	t possible	
d-squarks : kinematically not	t possible	
s-squarks : kinematically not	t possible	
b-squarks : kinematically not	t possible	
selectrons		

selectrons 1 1 290.9786253 fb

1 2	2 2	81.1156881 fb	
2	2	43.5380487 fb	
smuo	ons		
1	1	57.2730680 fb	
1	2	0.0058106 fb	
2	2	18.4801616 fb	
stau	ıs		
1	1	62.1659720 fb	
1	2	1.2151999 fb	
2	2	15.5585302 fb	
e-si	neuti	rino	
		412.6967832 fb	
mu-s	sneut	trino	
		12.9553932 fb	
tau	-sneı	utrino	
		13.2367076 fb	
	trali		
1	1	284.9858436 fb	
1	2	64.4974860 fb	
1	3	8.0550495 fb	
1	4	1.0888651 fb	
2	2	61.0982614 fb	
chai	rgind	09	
1	1	142.6642433 fb	
-	-		
h^0	Z		
		23.1467991 fb	
Н^О	z :	cross section below 1.0000000E	-03 fb
h^0	A^0	: kinematically not possible	
H ^ 0	A^0	: kinematically not possible	
H^+	H^-	: kinematically not possible	

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