Automated calculations for massive fermion production with "aTALC"

A. Lorca and T. Riemann

Deutsches Elektronen-Synchrotron, DESY, Platanenallee 6, 15738 Zeuthen, Germany

The package "aTALC" has been developed for the automated calculation of radiative corrections to two-fermion production at $e^+e^-$ colliders. The package uses DIANA, QGRAF, FORM, FORTRAN, FF, LOOPTOOLS, and further unix/linux tools. Numerical results are presented for $e^+e^- \rightarrow e^+e^-, \mu^+\mu^-, b\bar{b}, t\bar{t}, t\bar{c}$.

1. INTRODUCTION

Two fermion production, e.g.

$$e^+e^- \rightarrow e^+e^-, \mu^+\mu^-, \tau^+\tau^-, b\bar{b}, t\bar{t}, b\bar{s}, t\bar{c},$$

is among the reactions to be observed at a future linear $e^+e^-$ collider (LC) [1]. We developed packages for their calculation in the electroweak Standard Model (SM) (and extensions of it). The resulting FORTRAN codes may serve as etalons for other, numerical programs, may be used as electroweak library for some Monte Carlo program as well as directly used for studying the corresponding scattering process.

Earlier studies in this connection are [2–5]. Here, we report on "aTALC", a package with a high degree of automatization for this kind of calculations, and will focus on two applications: Bhabha scattering and flavour violating fermion pair production.

2. AUTOMATED CALCULATIONS WITH aTALC

The package "aTALC" for the automatic calculation of a variety of two fermion production processes may be obtained from [6] where also its installation will be described. The logical structure of the package is shown in Figure 1. It consists of three modules: DIANA, kitFORM3 and kitFORTRAN. In order to run and/or modify one of the samples in the example directory, the user has to choose the process in terms of incoming and outgoing fermions and the model lagrangian; we support two models: QED.model and EWSM.model, both with counterterms. This is done by modifying the driver file process.ini in the sample directory tree. Sample processes are: Bhabha scattering (eeee), $\mu$ pairs (muon production), $b$ pairs (eebb), $b\bar{s}$ (leLe-bS), $t\bar{c}$ (leLe-tC), etc. Selected cases will also be part of the public distribution.

The electroweak corrections are organized following [7,8], and we keep all the fermion masses, including $m_e$, by default.

The user runs the package by MAKE in e.g. eebb.user and produces (among others) the subdirectories tree, loop and fortran. The DIANA module uses QGRAF v.2.0 [9] and DIANA v.2.35 [10] and creates symbolic FORM-readable output for each of the Feynman diagrams in eebb.in. Moreover, graphical representations are produced, both stored as encapsulated postscript figures in EPS (single Feynman diagrams) and as postscripts in eebbInfo.ps (with detailed informations for each single diagram) and in eebb.ps (an overview of all contributions).

Then module kitFORM3 performs some algebraic simplifications and determines the matrix elements and form factors. This operation takes place separately for the tree and loop levels. The module kitFORTRAN returns the FORTRAN executable file main.out and the library libaitalc_v1.a. Two FORTRAN files, parameterlist.hf and main.f, provide the user access to the input parameters in the model.

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and to the design of numerical output of the code (i.e., number of data points in the angular distribution, integrated cross section, running flags, etc.). By executing `main.out`, the user produces a sample output `main.log` with differential and integrated cross sections, as well as forward-backward asymmetries. For Bhabha scattering at $\sqrt{s} = 500$ GeV, a shortened sample is reproduced in Table 1.

Advanced features are not extremely user friendly, but are still under development. Soft photonic corrections may be included (or not), and with `lidentCKM=.true`. quark flavour mixing is discarded with a diagonal CKM matrix. One may also perform a full one-loop tensor integral reduction to the master integrals $A_0$, $B_0$, $C_0$ and $D_0$ in the Passarino-Veltman scheme \[^{[11]}\]. All these possibilities will be described in more detail in a tutorial to be published soon; see also \[^{[12]}\].

3. SELECTED APPLICATIONS

3.1. Bhabha scattering

Bhabha scattering, 

$$e^+e^- \rightarrow e^+e^-,$$  \(^{(2)}\)

was one of the first processes to be calculated in QED \[^{[13]}\], and the one-loop corrections in the SM were treated in \[^{[14–20]}\].

We are calculating the virtual corrections to massive, low angle Bhabha scattering at a Linear Collider \[^{[1]}\]. In order to guarantee a precision of $10^{-4}$ of the differential cross section we need the one-loop corrections in the SM, and the two-loop corrections in pure, massive QED \[^{[21]}\]; the latter comprise also the squared one-loop corrections \[^{[22]}\].
The one-loop SM corrections are determined with \texttt{aTALC}. A shortened example of the \texttt{aTALC} output is given in Table \ref{table:output}. A comparison of our calculation with another one based on \texttt{FORMCALC} \cite{23} is reproduced in \cite{12}, where an agreement is obtained of more than 12 significant digits.

3.2. At the $Z$-peak

For the numerical evaluation of the one-loop functions we use \texttt{LoopTools} v.2.0 and follow the conventions given in the manual \cite{24–26}. \texttt{LoopTools} relies on the package FF \cite{25}. In contrast to FF, it does not support the case of complex masses. Near the $Z$-peak, the Breit-Wigner propagator has to be used,

\[
\frac{1}{s - m_Z^2 + i\varepsilon} \rightarrow \frac{1}{s - m_Z^2 + i\Gamma_Z m_Z},
\]

with a width parameter $\Gamma_Z$. Flag \texttt{width} activates the following changes in \texttt{aTALC}:

- Apply substitution \texttt{K} in Feynman rules;
- Avoid double counting by discarding the $Z$ self-energy diagrams resummed in $\Gamma_Z$;
- Use loop integrals with complex $Z$ mass.

In general we use for the numerics Passarino-Veltman functions of scalar, vector, and tensor type. In the neighbourhood of the $Z$ resonance, though, we perform the tensor reduction for the $\gamma - Z$ dependent functions $D_{ij}(...)$, and have to calculate the following functions with complex $Z$ mass: $B_0$, $C_0$ and $D_0$. The $A_0$ and $B_0$ are trivial, and the infrared divergent $D_0$ was calculated in \cite{27}. The tensor reduction of the tensor integrals related to the $\gamma - Z$ box diagrams leads, by shrinking of internal lines, to infrared safe two- and three-point functions with complex $Z$ mass. They are generally treated in \cite{28} and available in the FF package \cite{25} \footnote{It has to be noted that the packages FF and \texttt{LoopTools} v.2.0 \cite{24} are not compatible and the user has to either drop one of them or modify the internal source code. Therefore, presently \texttt{aTALC} renames part of the \texttt{LoopTools} subroutines. After this workshop, \texttt{LoopTools} v.2.1 (29 June 2004) was released and is now prepared for the treatment of complex masses.}.

\begin{equation}
\sigma_{b\bar{s}}(\sqrt{s} = m_Z) = (1.136 \pm 0.001) \text{ fb},
\end{equation}

\begin{equation}
\sigma_{b\bar{s}}(\sqrt{s} = 200\text{GeV}) = (2.033 \pm 0.002) \text{ fb}.
\end{equation}

For the other channel, we agree within the accuracy of the figures. Figure \ref{fig:total} shows also the total cross section for $b\bar{s}$.

3.3. Flavour number violation

Topics of particular interest are flavour changing neutral current processes at $e^+e^-$ colliders since they are forbidden in the Born approximation of the Standard Model and might indicate New Physics. At one-loop order, they may occur in the SM if the fermions of different flavour have different masses and are mixing. Usually one calculates simply the flavour changing $Z$ decay rate because the chances to observe an effect are largest at the $Z$-peak. The predictions in the minimally extended Standard Model are tiny (see e.g. \cite{29–31}), but may be much larger with supersymmetry (see e.g. \cite{32}). Recently, there were also studies of the complete scattering processes \cite{33,34}:

\[
e^+e^- \rightarrow \mu^+\mu^- \text{ around the $Z$-peak is presented in Figure \ref{fig:muon}.}
\]

4. CONCLUDING REMARKS

We plan to include into \texttt{aTALC} also processes which have contributions from Feynman diagrams with five-point functions. This would allow us to
Table 1
Sample output of aITALC for Bhabha scattering at $\sqrt{s} = 500$ GeV: Born and 1-loop corrected differential cross sections; input data as in [3,12]

```plaintext
# aITALC: Version 0.7 by A.Lorca -- T.Riemann
# ==================================================================
#cos(theta)  dcs(BORN) ... dcs(BORN+Q+W+soft) ...
-0.90000  0.2169988288109205E+00 ... 0.1934450785268578E+00 ...
-0.50000  0.2613604305853236E+00 ... 0.2387066977233451E+00 ...
 0.00000  0.5981423072503301E+00 ... 0.5466771794694227E+00 ...
 0.50000  0.4212729493916255E+01 ... 0.3813007881789546E+01 ...
 0.90000  0.1891603223322704E+03 ... 0.1729283490665079E+03 ...
# ==================================================================
```

Figure 2. The infrared finite function $C_0(s, m_{\mu}^2, m_{\mu}^2, m_Z^2, 0, m_{\mu}^2)$ as a function of $\sqrt{s}$.

Figure 3. Muon production at the $Z$-peak.

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Differential Cross Section

Figure 4. Differential cross sections for $e^+e^- \rightarrow b\bar{s}$ and $e^+e^- \rightarrow t\bar{c}$ at $\sqrt{s} = 200$ GeV; input data as in [34].

Total Cross Section: $e^+e^- \rightarrow b\bar{s}$

Figure 5. Integrated cross section for $e^+e^- \rightarrow b\bar{s}$ as a function of $\sqrt{s}$; input data as in [33].