

SMEFTsim 3.0 — a practical guide

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ABSTRACT: The `SMEFTsim` package [1] is designed to enable automated computations in the Standard Model Effective Field Theory (SMEFT), where the SM Lagrangian is extended with a complete basis of dimension six operators. It contains a set of models written in `FeynRules` and pre-exported to the `UFO` format, for usage within Monte Carlo event generators. The models differ in the flavor assumptions and in the input parameters chosen for the electroweak sector. The present document provides a self-contained, pedagogical reference that collects all the theoretical and technical aspects relevant to the use of `SMEFTsim` and it documents the release of version 3.0. Compared to the previous release, the description of Higgs production via gluon-fusion in the SM has been significantly improved, two flavor assumptions for studies in the top quark sector have been added, and a new feature has been implemented, that enables the treatment of linearized SMEFT corrections to the propagators of unstable particles.

`SMEFTsim` 3.0 is available on the Github website <https://SMEFTsim.github.io> and on the `FeynRules` database <http://feynrules.irmp.ucl.ac.be/wiki/SMEFT>.

KEYWORDS: Beyond Standard Model, Effective Field Theories, Higgs Physics

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1 Introduction

LHC physics is about to enter a precision era that will span over the next two decades. During this time, new opportunities to hunt for new physics will arise: direct searches of new particles will be complemented by indirect searches, that target possible deviations from the predictions of the Standard Model (SM). While the isolation of this kind of signatures is not without challenges, indirect searches present some very attractive features. Most notably, they do not rely on specific assumptions about the nature of the new physics under scrutiny and, at the same time, their sensitivity in terms of new physics scales can potentially extend beyond the energy reach of the collider.

The Standard Model Effective Field Theory (SMEFT) is the best established theory framework to describe such effects. Its formulation employs the degrees of freedom and gauge symmetries of the SM and it is structured as an infinite series of operators sorted by canonical dimension. At the observables level, it reproduces a series expansion in (E/Λ) , being E the typical energy exchanged in a process and Λ the mass scale that characterizes the beyond-SM (BSM) dynamics. The condition $(E/\Lambda) \ll 1$, indicating the near decoupling of the new physics sector, is necessarily assumed.

The SMEFT has been developed extensively in the past ten years, laying the ground for a systematic program for indirect searches [2–4]. The ultimate goal is to measure as many EFT parameters as possible, in a manner that enables the extraction of unbiased information about the underlying physics. The crucial aspect of this program is its transversality: the SMEFT contains a large number of parameters, each typically entering the description

of several processes. Combining measurements of different observables is then mandatory in order to preserve the model-independence of the analysis. To date, this principle has been applied within individual sectors as well as across Higgs, electroweak (EW) and top quark measurements, see refs. [5–15] for recent examples. The incorporation of data from flavor observables (including non-LHC experiments) would be very valuable in this context, as most of the SMEFT parameter space is “flavorful”. First steps in this direction were taken in [16–19].

The theory developments have been accompanied by the publication of a number of computing tools that automate most stages of a SMEFT study [20]. These include the definition of non-redundant operator bases and the translation between them [21–24], the matching to concrete BSM models or to the low-energy EFT and the renormalization group running [25–31], the extraction of the Feynman rules in R_ξ gauges [32, 33] and in the background field gauge [34], Monte Carlo simulations [1, 35–37] and global analyses [13, 38–41].

The `SMEFTsim` package [1] was designed in order to enable the Monte Carlo simulation of arbitrary processes in the effective theory, in the spirit of providing a unified, general-purpose tool for SMEFT physics at the LHC. It provides complete tree level, unitary gauge predictions at $\mathcal{O}(\Lambda^{-2})$, including *all* the dimension six operators in the so-called Warsaw basis [42]. The field and parameter redefinitions that are required in order to compute physical observables in the SMEFT are conveniently performed internally. The package contains `FeynRules` [43, 44] source files and a set of models pre-exported to the `UFO` format [45]. Although the latter are in principle compatible with most Monte Carlo generators, they have been optimized for the use in `MadGraph5_aMC@NLO` [46], that is illustrated in section 8. Potential issues due, for instance, to the event generation entering regions where the EFT validity or the unitarity of the S -matrix are violated are not addressed within `SMEFTsim` itself, but can be generally treated with tools offered by the Monte Carlo generators or with theoretical assessments *a posteriori*.

The `UFO` models differ in the flavor assumptions and in the choice of the input parameters for the EW sector.¹ The original release implemented three alternative flavor scenarios: a general one, a $U(3)^5$ -symmetric case and a linear minimal flavor violation (MFV) option where BSM CP-violating phases are forbidden. For each setup, it offered two EW input sets: $\{\alpha_{\text{em}}, m_Z, G_F\}$ or $\{m_W, m_Z, G_F\}$.

This work documents the release of `SMEFTsim` version 3.0, that introduces a number of improvements summarized below. It is also meant as a pedagogical and self-contained reference for its usage, where all the relevant theoretical aspects are reviewed in detail.

The present document is structured as follows: sections 1–4 review the theoretical background while sections 5–8 describe technical aspects of the `FeynRules` and `UFO` implementations and provide recommendations for their use.

The notation is fixed in section 1.1. Section 2 focuses on the bosonic sector and it reviews the field and parameter redefinitions required to ensure a canonical parameterization of the kinetic terms and scalar potential. Section 3 is devoted to the flavor structure of

¹The original release contained two fully equivalent implementations, that were called model sets A and B. Both were provided for debugging and cross-validation. Set B is not supported anymore starting from version 3.0, which is based on set A.

the SMEFT and it defines the five scenarios implemented in `SMEFTsim`. Some significant changes have been made compared to version 2, that are documented in detail, and two new flavor options have been introduced (`top`, `topU31`) that comply with the recommendations for studies of top quark observables [36]. Section 4 provides a general discussion of how the extraction of numerical values for the SM parameters is affected in the presence of higher-dimensional operators, and illustrates the treatment of these effects in `SMEFTsim`.

Section 5 documents the implementation of Higgs interactions that are purely loop-generated in the SM, namely $h\gamma\gamma$, $hZ\gamma$, hgg : as `SMEFTsim` only supports tree-level interactions, these are treated as effective vertices in the large top mass limit, which is a good approximation for Higgs production and decay processes. Compared to version 2, the description of Higgs-gluon vertices has been substantially improved, such that it can now model one-loop SM interactions with up to 4 gluons. Section 6 focuses on SMEFT effects in the propagators of unstable particles, that arise due to modifications of their pole masses or decay widths. A new feature has been introduced in version 3.0, that enables the inclusion of such effects, linearized in the EFT parameters, in Monte Carlo simulations. To our knowledge `SMEFTsim` is the first publicly available UFO model to implement such a tool. Sections 7, 8 provide recommendations for the use of `SMEFTsim` in `Mathematica` and in `MadGraph5_aMC@NLO` respectively, and in section 9 we conclude.

Additional useful material is provided in the appendices: analytic expressions of the decay widths implemented in the propagator corrections (appendix A), a list of changes made in version 3.0 (appendix B), tables to facilitate the conversion between flavor assumptions (appendix C), between theory and code notation (appendix D) and between `SMEFTsim` and `dim6top` or `SMEFT@NLO` (appendix E). Finally, appendix F documents the validation of the UFO models, that followed the procedure recommended in [47].

1.1 Basics and notation

We consider the SMEFT Lagrangian truncated at the dimension-6 level:

$$\mathcal{L}_{\text{SMEFT}} = \mathcal{L}_{\text{SM}} + \mathcal{L}_6. \tag{1.1}$$

We neglect all lepton- and baryon-number violating terms, which includes the dimension-5 Weinberg operator that generates a Majorana mass term for neutrinos. For future convenience, the SM Lagrangian is split into four terms:

$$\mathcal{L}_{\text{SM}} = \mathcal{L}_{\text{gauge}} + \mathcal{L}_{\text{fermions}} + \mathcal{L}_{\text{Yukawa}} + \mathcal{L}_{\text{Higgs}}, \tag{1.2}$$

where

$$\mathcal{L}_{\text{gauge}} = -\frac{1}{4}B_{\mu\nu}B^{\mu\nu} - \frac{1}{4}W_{\mu\nu}^i W^{i\mu\nu} - \frac{1}{4}G_{\mu\nu}^a G^{a\mu\nu}, \tag{1.3}$$

$$\mathcal{L}_{\text{fermions}} = \bar{q}i\not{D}q + \bar{u}i\not{D}u + \bar{d}i\not{D}d + \bar{l}i\not{D}l + \bar{e}i\not{D}e, \tag{1.4}$$

$$\mathcal{L}_{\text{Yukawa}} = -\bar{d}Y_d H^\dagger q - \bar{u}Y_u \tilde{H}^\dagger q - \bar{e}Y_l H^\dagger l + \text{h.c.}, \tag{1.5}$$

$$\mathcal{L}_{\text{Higgs}} = D_\mu H^\dagger D^\mu H + m^2(H^\dagger H) - \lambda(H^\dagger H)^2. \tag{1.6}$$

q, l represent the left-handed quark and lepton doublets respectively, and u, d, e the right-handed quarks and leptons. H is the Higgs doublet and $\tilde{H} = i\sigma^2 H^*$, where $\sigma^i, i = \{1, 2, 3\}$ are the Pauli matrices. Y_d, Y_u, Y_l are the 3×3 Yukawa matrices of the down and up quarks and of the charged leptons. Covariant derivatives are defined with a plus sign, i.e.²

$$D_\mu q = \left[\partial_\mu + ig_s T^a G_\mu^a + i \frac{g_W}{2} \sigma^i W_\mu^i + i \mathbf{y}_q g_1 B_\mu \right] q. \quad (1.7)$$

$T^a \equiv \lambda^a/2, a = \{1, \dots, 8\}$ are the $SU(3)_c$ generators, with λ^a the Gell-Mann matrices. $\mathbf{y}_q = 1/6$ denotes the hypercharge of the q field and g_s, g_W, g_1 are the $SU(3)_c \times SU(2)_L \times U(1)_Y$ coupling constants. As a general rule, color indices are denoted by a, b, c, d , $SU(2)_L$ indices by i, j, k and flavor indices by p, r, s, t . Summation over identical indices is always understood, unless otherwise specified.

The Lagrangian \mathcal{L}_6 contains a complete and non-redundant basis of dimension-6 operators Q_α constructed with the SM fields and invariant under the $SU(3)_c \times U(2)_L \times U(1)_Y$ gauge symmetry. SMEFTsim implements the Warsaw basis [42], whose operators are collected in 8 groups, following the classification of ref. [48]. Class 8 is further split into 4 subgroups.³

$$\mathcal{L}_6 = \mathcal{L}_6^{(1)} + \mathcal{L}_6^{(2)} + \mathcal{L}_6^{(3)} + \mathcal{L}_6^{(4)} + \mathcal{L}_6^{(7)} + \mathcal{L}_6^{(8)} + \left[\mathcal{L}_6^{(5)} + \mathcal{L}_6^{(6)} + \text{h.c.} \right], \quad (1.8)$$

$$\mathcal{L}_6^{(8)} = \mathcal{L}_6^{(8a)} + \mathcal{L}_6^{(8b)} + \mathcal{L}_6^{(8c)} + \left[\mathcal{L}_6^{(8d)} + \text{h.c.} \right]. \quad (1.9)$$

Each sub-Lagrangian has the form

$$\mathcal{L}_6^{(n)} = \frac{1}{\Lambda^2} \sum_\alpha C_\alpha Q_\alpha, \quad (1.10)$$

with the sum running over the class- n operators $\{Q_\alpha\}$ defined in table 1 and C_α denoting the associated Wilson coefficients. Both Q_α and C_α generally carry flavor indices, that are implicitly contracted in eq. (1.10). In this basis, explicit CP violation is carried by the real coefficients $C_{\tilde{G}}, C_{\tilde{W}}, C_{H\tilde{G}}, C_{H\tilde{W}}, C_{H\tilde{B}}, C_{H\tilde{W}B}$ and by the imaginary parts of the Wilson coefficients associated to non-Hermitian fermionic operators, namely those in $\mathcal{L}_6^{(5),(6),(8d)}$ and \mathcal{O}_{Hud} . Baryon-number violating operators are omitted.

The operators definitions use the following notation:

$$\tilde{X}^{\mu\nu} = \frac{1}{2} \varepsilon^{\mu\nu\rho\sigma} X_{\rho\sigma}, \quad H^\dagger i \overleftrightarrow{D}_\mu H = H^\dagger (iD_\mu H) - (iD_\mu H^\dagger) H, \quad (1.11)$$

$$\sigma^{\mu\nu} = \frac{i}{2} [\gamma^\mu, \gamma^\nu], \quad H^\dagger i \overleftrightarrow{D}_\mu^i H = H^\dagger \sigma^i (iD_\mu H) - (iD_\mu H^\dagger) \sigma^i H. \quad (1.12)$$

²The covariant derivative sign is handled automatically by **FeynRules**. The convention chosen here also implies that gauge field strengths have the form $W_{\mu\nu}^i = \partial_\mu W_\nu^i - \partial_\nu W_\mu^i - g_W \varepsilon^{ijk} W_\mu^j W_\nu^k$, etc.

³Note that $\mathcal{L}_6^{(7)}$ implicitly contains $(Q_{Hud} + \text{h.c.})$, as this operator is not Hermitian.

$\mathcal{L}_6^{(1)} - X^3$		$\mathcal{L}_6^{(6)} - \psi^2 XH$		$\mathcal{L}_6^{(8b)} - (\bar{R}R)(\bar{R}R)$	
Q_G	$f^{abc}G_\mu^{a\nu}G_\nu^{b\rho}G_\rho^{c\mu}$	Q_{eW}	$(\bar{l}_p\sigma^{\mu\nu}e_r)\sigma^iHW_\mu^i$	Q_{ee}	$(\bar{e}_p\gamma_\mu e_r)(\bar{e}_s\gamma^\mu e_t)$
$Q_{\tilde{G}}$	$f^{abc}\tilde{G}_\mu^{a\nu}G_\nu^{b\rho}G_\rho^{c\mu}$	Q_{eB}	$(\bar{l}_p\sigma^{\mu\nu}e_r)HB_{\mu\nu}$	Q_{uu}	$(\bar{u}_p\gamma_\mu u_r)(\bar{u}_s\gamma^\mu u_t)$
Q_W	$\varepsilon^{ijk}W_\mu^{i\nu}W_\nu^{j\rho}W_\rho^{k\mu}$	Q_{uG}	$(\bar{q}_p\sigma^{\mu\nu}T^a u_r)\tilde{H}G_\mu^a$	Q_{dd}	$(\bar{d}_p\gamma_\mu d_r)(\bar{d}_s\gamma^\mu d_t)$
$Q_{\tilde{W}}$	$\varepsilon^{ijk}\tilde{W}_\mu^{i\nu}W_\nu^{j\rho}W_\rho^{k\mu}$	Q_{uW}	$(\bar{q}_p\sigma^{\mu\nu}u_r)\sigma^i\tilde{H}W_\mu^i$	Q_{eu}	$(\bar{e}_p\gamma_\mu e_r)(\bar{u}_s\gamma^\mu u_t)$
$\mathcal{L}_6^{(2)} - H^6$		Q_{uB}	$(\bar{q}_p\sigma^{\mu\nu}u_r)\tilde{H}B_{\mu\nu}$	Q_{ed}	$(\bar{e}_p\gamma_\mu e_r)(\bar{d}_s\gamma^\mu d_t)$
Q_H	$(H^\dagger H)^3$	Q_{dG}	$(\bar{q}_p\sigma^{\mu\nu}T^a d_r)HG_\mu^a$	$Q_{ud}^{(1)}$	$(\bar{u}_p\gamma_\mu u_r)(\bar{d}_s\gamma^\mu d_t)$
$\mathcal{L}_6^{(3)} - H^4 D^2$		Q_{dW}	$(\bar{q}_p\sigma^{\mu\nu}d_r)\sigma^iHW_\mu^i$	$Q_{ud}^{(8)}$	$(\bar{u}_p\gamma_\mu T^a u_r)(\bar{d}_s\gamma^\mu T^a d_t)$
$Q_{H\Box}$	$(H^\dagger H)\Box(H^\dagger H)$	Q_{dB}	$(\bar{q}_p\sigma^{\mu\nu}d_r)HB_{\mu\nu}$		
Q_{HD}	$(D^\mu H^\dagger H)(H^\dagger D_\mu H)$				
$\mathcal{L}_6^{(4)} - X^2 H^2$		$\mathcal{L}_6^{(7)} - \psi^2 H^2 D$		$\mathcal{L}_6^{(8c)} - (\bar{L}L)(\bar{R}R)$	
Q_{HG}	$H^\dagger HG_\mu^a G^{a\mu\nu}$	$Q_{Hl}^{(1)}$	$(H^\dagger i\overleftrightarrow{D}_\mu H)(\bar{l}_p\gamma^\mu l_r)$	Q_{le}	$(\bar{l}_p\gamma_\mu l_r)(\bar{e}_s\gamma^\mu e_t)$
$Q_{H\tilde{G}}$	$H^\dagger H\tilde{G}_\mu^a G^{a\mu\nu}$	$Q_{Hl}^{(3)}$	$(H^\dagger i\overleftrightarrow{D}_\mu^i H)(\bar{l}_p\sigma^i\gamma^\mu l_r)$	Q_{lu}	$(\bar{l}_p\gamma_\mu l_r)(\bar{u}_s\gamma^\mu u_t)$
Q_{HW}	$H^\dagger HW_\mu^i W^{I\mu\nu}$	Q_{He}	$(H^\dagger i\overleftrightarrow{D}_\mu H)(\bar{e}_p\gamma^\mu e_r)$	Q_{ld}	$(\bar{l}_p\gamma_\mu l_r)(\bar{d}_s\gamma^\mu d_t)$
$Q_{H\tilde{W}}$	$H^\dagger H\tilde{W}_\mu^i W^{i\mu\nu}$	$Q_{Hq}^{(1)}$	$(H^\dagger i\overleftrightarrow{D}_\mu H)(\bar{q}_p\gamma^\mu q_r)$	Q_{qe}	$(\bar{q}_p\gamma_\mu q_r)(\bar{e}_s\gamma^\mu e_t)$
Q_{HB}	$H^\dagger HB_{\mu\nu}B^{\mu\nu}$	$Q_{Hq}^{(3)}$	$(H^\dagger i\overleftrightarrow{D}_\mu^i H)(\bar{q}_p\sigma^i\gamma^\mu q_r)$	$Q_{qu}^{(1)}$	$(\bar{q}_p\gamma_\mu q_r)(\bar{u}_s\gamma^\mu u_t)$
$Q_{H\tilde{B}}$	$H^\dagger H\tilde{B}_{\mu\nu}B^{\mu\nu}$	Q_{Hu}	$(H^\dagger i\overleftrightarrow{D}_\mu H)(\bar{u}_p\gamma^\mu u_r)$	$Q_{qu}^{(8)}$	$(\bar{q}_p\gamma_\mu T^a q_r)(\bar{u}_s\gamma^\mu T^a u_t)$
Q_{HWB}	$H^\dagger\sigma^i HW_\mu^i B^{\mu\nu}$	Q_{Hd}	$(H^\dagger i\overleftrightarrow{D}_\mu H)(\bar{d}_p\gamma^\mu d_r)$	$Q_{qd}^{(1)}$	$(\bar{q}_p\gamma_\mu q_r)(\bar{d}_s\gamma^\mu d_t)$
$Q_{H\tilde{W}B}$	$H^\dagger\sigma^i H\tilde{W}_\mu^i B^{\mu\nu}$	$Q_{Hud} + \text{h.c.}$	$i(\tilde{H}^\dagger D_\mu H)(\bar{u}_p\gamma^\mu d_r)$	$Q_{qd}^{(8)}$	$(\bar{q}_p\gamma_\mu T^a q_r)(\bar{d}_s\gamma^\mu T^a d_t)$
$\mathcal{L}_6^{(5)} - \psi^2 H^3$		$\mathcal{L}_6^{(8a)} - (\bar{L}L)(\bar{L}L)$		$\mathcal{L}_6^{(8d)} - (\bar{L}R)(\bar{R}L), (\bar{L}R)(\bar{L}R)$	
Q_{eH}	$(H^\dagger H)(\bar{l}_p e_r H)$	Q_{ll}	$(\bar{l}_p\gamma_\mu l_r)(\bar{l}_s\gamma^\mu l_t)$	Q_{ledq}	$(\bar{l}_p^j e_r)(\bar{d}_s q_t^j)$
Q_{uH}	$(H^\dagger H)(\bar{q}_p u_r \tilde{H})$	$Q_{qq}^{(1)}$	$(\bar{q}_p\gamma_\mu q_r)(\bar{q}_s\gamma^\mu q_t)$	$Q_{quqd}^{(1)}$	$(\bar{q}_p^j u_r)\varepsilon_{jk}(\bar{q}_s^k d_t)$
Q_{dH}	$(H^\dagger H)(\bar{q}_p d_r H)$	$Q_{qq}^{(3)}$	$(\bar{q}_p\gamma_\mu\sigma^i q_r)(\bar{q}_s\gamma^\mu\sigma^i q_t)$	$Q_{quqd}^{(8)}$	$(\bar{q}_p^j T^a u_r)\varepsilon_{jk}(\bar{q}_s^k T^a d_t)$
		$Q_{lq}^{(1)}$	$(\bar{l}_p\gamma_\mu l_r)(\bar{q}_s\gamma^\mu q_t)$	$Q_{lequ}^{(1)}$	$(\bar{l}_p^j e_r)\varepsilon_{jk}(\bar{q}_s^k u_t)$
		$Q_{lq}^{(3)}$	$(\bar{l}_p\gamma_\mu\sigma^i l_r)(\bar{q}_s\gamma^\mu\sigma^i q_t)$	$Q_{lequ}^{(3)}$	$(\bar{l}_p^j\sigma_{\mu\nu}e_r)\varepsilon_{jk}(\bar{q}_s^k\sigma^{\mu\nu}u_t)$

Table 1. \mathcal{L}_6 operators in the Warsaw basis [42], categorized into eight classes $\mathcal{L}_6^{(n)}$ as in [48]. Only baryon number-conserving invariants are retained. The flavor indices p, r, s, t are suppressed in the operators' names.

2 EWSB, field and parameter redefinitions

This section reviews the Lagrangian manipulations that are required in order to compute physical processes in the SMEFT truncated at $\mathcal{O}(\Lambda^{-2})$.⁴The procedure described here largely overlaps with what reported e.g. in [1, 32, 48–59].

2.1 Higgs sector

The operator Q_H introduces a perturbation of the Higgs potential:

$$V(H) = -m^2 H^\dagger H + \lambda(H^\dagger H)^2 - \frac{C_H}{\Lambda^2}(H^\dagger H)^3. \quad (2.1)$$

The true minimum of the potential, that triggers the electroweak symmetry breaking, is

$$\begin{aligned} \langle H^\dagger H \rangle &= \frac{v_T^2}{2} \equiv \frac{v^2}{2} \left[1 + \frac{3}{4\lambda} \frac{v^2}{\Lambda^2} C_H \right] + \mathcal{O}(\Lambda^{-4}) \\ &= \frac{v^2}{2} \left[1 + \frac{3}{4\lambda} \bar{C}_H \right] + \mathcal{O}(\Lambda^{-4}), \end{aligned} \quad (2.2)$$

with

$$v^2 = \frac{m^2}{2\lambda}. \quad (2.3)$$

We have introduced the “bar” notation for Wilson coefficients:

$$\bar{C}_\alpha \equiv \frac{v_T^2}{\Lambda^2} C_\alpha. \quad (2.4)$$

Note that because $v = v_T + \mathcal{O}(\Lambda^{-2})$ and $\mathcal{O}(\Lambda^{-4})$ contributions are entirely neglected, the two quantities v, v_T are interchangeable whenever they multiply a Wilson coefficient.⁵

The Higgs field H is expanded around its vacuum expectation value (vev) as

$$H = \begin{pmatrix} -iG^+ \\ \frac{v_T + h + iG^0}{\sqrt{2}} \end{pmatrix}, \quad (2.5)$$

with G^+, G^0 the charged and neutral Goldstone bosons and h the physical Higgs boson. In the broken phase, the kinetic terms of the scalar fields receive corrections from the operators $Q_{H\Box}, Q_{HD}$. As the scope of `SMEFTsim` is limited to tree-level calculations, we choose to work in unitary gauge and neglect EFT effects in the Goldstone sector, both in the present discussion and in the code implementations. The Goldstone bosons case and the generalization of the gauge fixing procedure in the SMEFT were addressed in refs. [3, 32, 58–62].

⁴Within a Monte Carlo event generation, `SMEFTsim` generally enables the computation of higher order corrections to a given observable, such as $\mathcal{O}(\Lambda^{-4})$ corrections stemming from the square of $\mathcal{O}(\Lambda^{-2})$ amplitudes (see section 8). However, consistent results are only provided to $\mathcal{O}(\Lambda^{-2})$, as the SMEFT Lagrangian implemented is truncated at this order.

⁵In fact, it would be more appropriate to define the \bar{C}_α notation with the parameter \hat{v} defined in section 4.2, rather than with v_T . However, as long as $\mathcal{O}(\Lambda^{-4})$ terms are neglected, all three v 's are formally identical when multiplying a Wilson coefficient.

Using integration by parts, the kinetic term of the physical Higgs boson takes the form

$$\mathcal{L}_{\text{Higgs}} + \mathcal{L}_6 = \frac{1}{2} \partial_\mu h \partial^\mu h [1 - 2\Delta\kappa_H] + \dots \quad \text{with} \quad \Delta\kappa_H = \bar{C}_{H\Box} - \frac{\bar{C}_{HD}}{4}, \quad (2.6)$$

and it is brought to the canonical normalization via the field redefinition

$$h \rightarrow [1 + \Delta\kappa_H] h. \quad (2.7)$$

This replacement is formally operated on the entire $\mathcal{L}_{\text{SMEFT}}$. However, when applied to \mathcal{L}_6 , its net effect is of $\mathcal{O}(\Lambda^{-4})$. As we work at $\mathcal{O}(\Lambda^{-2})$, the replacement only needs to be performed on \mathcal{L}_{SM} . This holds for all field and parameter redefinitions introduced in the following, unless otherwise specified. For the same reason, all quantities in a Wilson coefficient's prefactor are understood to be defined in the SM limit.

The main consequence of (2.7) is that the Wilson coefficients $C_{H\Box}$, C_{HD} are recast into an overall rescaling of all SM Higgs couplings. The resulting Higgs potential is

$$\begin{aligned} V(H) = & h^2 \lambda v_T^2 \left[1 + 2\Delta\kappa_H - \frac{3}{2\lambda} \bar{C}_H \right] + h^3 \lambda v_T \left[1 + 3\Delta\kappa_H - \frac{5}{2\lambda} \bar{C}_H \right] \\ & + h^4 \frac{\lambda}{4} \left[1 + 4\Delta\kappa_H - \frac{15}{2\lambda} \bar{C}_H \right] - \frac{3}{4} \frac{h^5}{v_T} \bar{C}_H - \frac{1}{8} \frac{h^6}{v_T^2} \bar{C}_H. \end{aligned} \quad (2.8)$$

In the `FeynRules` implementation, the redefinitions of the physical Higgs field, eq. (2.7), and of the vev, eq. (2.2), are embedded in the definition of the Higgs doublet.

2.2 Gauge sector

Upon EWSB, the operators Q_{HG} , Q_{HW} , Q_{HB} , Q_{HWB} induce corrections to the kinetic terms of the gauge bosons. The first three lead to overall rescalings:

$$\mathcal{L}_{\text{gauge}} + \mathcal{L}_6 = -\frac{1}{4} B_{\mu\nu} B^{\mu\nu} [1 - 2\bar{C}_{HB}] - \frac{1}{4} W_{\mu\nu}^i W^{i\mu\nu} [1 - 2\bar{C}_{HW}] - \frac{1}{4} G_{\mu\nu}^a G^{a\mu\nu} [1 - 2\bar{C}_{HG}] + \dots \quad (2.9)$$

where the dots stand for all other interaction terms induced by \mathcal{L}_6 . The canonical normalization is easily restored at $\mathcal{O}(\Lambda^{-2})$, via the field redefinitions

$$G_\mu^a \rightarrow G_\mu^a (1 + \bar{C}_{HG}), \quad W_\mu^i \rightarrow W_\mu^i (1 + \bar{C}_{HW}), \quad B_\mu \rightarrow B_\mu (1 + \bar{C}_{HB}). \quad (2.10)$$

In order to leave the covariant derivatives unchanged, the coupling constants need to be redefined at the same time. Neglecting $\mathcal{O}(\Lambda^{-4})$ corrections:

$$g_s \rightarrow g_s (1 - \bar{C}_{HG}), \quad g_W \rightarrow g_W (1 - \bar{C}_{HW}), \quad g_1 \rightarrow g_1 (1 - \bar{C}_{HB}). \quad (2.11)$$

The operator Q_{HWB} introduces a kinetic mixing between the B and W^3 fields of the form

$$- \frac{C_{HWB}}{2} \frac{v_T^2}{\Lambda^2} W_{\mu\nu}^3 B^{\mu\nu}. \quad (2.12)$$

The rotation [48]

$$\begin{pmatrix} W_\mu^3 \\ B_\mu \end{pmatrix} \rightarrow \begin{pmatrix} 1 & -\bar{C}_{HWB}/2 \\ -\bar{C}_{HWB}/2 & 1 \end{pmatrix} \begin{pmatrix} W_\mu^3 \\ B_\mu \end{pmatrix}, \quad (2.13)$$

removes this residual mixing and leads to fully canonical and diagonal kinetic terms. Once eqs. (2.10), (2.11), (2.13) have been applied, the electric-charge eigenstates W^\pm are obtained via the usual rotation

$$\begin{pmatrix} W_\mu^1 \\ W_\mu^2 \end{pmatrix} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ i & -i \end{pmatrix} \begin{pmatrix} W_\mu^+ \\ W_\mu^- \end{pmatrix}, \quad (2.14)$$

while the mass matrix of the neutral bosons is diagonalized by

$$\begin{pmatrix} W_\mu^3 \\ B_\mu \end{pmatrix} = \frac{1}{\sqrt{g_1^2 + g_W^2}} \begin{pmatrix} g_W & g_1 \\ -g_1 & g_W \end{pmatrix} \begin{pmatrix} 1 & \frac{1}{2} \frac{g_W^2 - g_1^2}{g_W^2 + g_1^2} \bar{C}_{HWB} \\ -\frac{1}{2} \frac{g_W^2 - g_1^2}{g_W^2 + g_1^2} \bar{C}_{HWB} & 1 \end{pmatrix} \begin{pmatrix} Z_\mu \\ A_\mu \end{pmatrix}, \quad (2.15)$$

The rightmost rotation is unitary up to $\mathcal{O}(\Lambda^{-4})$ corrections, and therefore does not reintroduce kinetic mixing at $d = 6$. Equivalently,

$$\begin{pmatrix} W_\mu^3 \\ B_\mu \end{pmatrix} = \begin{pmatrix} \cos \theta & \sin \theta \\ -\sin \theta & \cos \theta \end{pmatrix} \begin{pmatrix} Z_\mu \\ A_\mu \end{pmatrix}, \quad (2.16)$$

with a shifted weak mixing angle θ defined as

$$\tan \theta = \frac{g_1}{g_W} + \frac{1}{2} \bar{C}_{HWB} \left(1 - \frac{g_1^2}{g_W^2} \right). \quad (2.17)$$

After all the coupling and field redefinitions have been applied, a generic covariant derivative has the form

$$D_\mu = \partial_\mu + iQ \frac{g_1 g_W}{\sqrt{g_1^2 + g_W^2}} A_\mu \left[1 - \bar{C}_{HWB} \frac{g_1 g_W}{g_W^2 + g_1^2} \right] \quad (2.18)$$

$$+ i\sqrt{g_1^2 + g_W^2} Z_\mu \left[T_3 - \frac{g_1^2}{g_1^2 + g_W^2} Q + \bar{C}_{HWB} \frac{g_1 g_W}{g_1^2 + g_W^2} \left(T_3 - \frac{g_W^2}{g_1^2 + g_W^2} Q \right) \right]$$

+ ...

$$= \partial_\mu + iQ g_W s_\theta A_\mu \left[1 - \frac{1}{2} \frac{c_\theta}{s_\theta} \bar{C}_{HWB} \right] + i \frac{g_W}{c_\theta} Z_\mu (T_3 - Q s_\theta^2) \left[1 + \frac{1}{2} \frac{s_\theta}{c_\theta} \bar{C}_{HWB} \right] \quad (2.19)$$

+ ...

where T_3 denotes the eigenvalue of the 3rd $SU(2)_L$ generator ($T_3 = \pm 1/2$ for left-handed fields and $T_3 = 0$ for right-handed ones) and $Q = T_3 + \mathbf{y}$ is the electric charge. We have also introduced the shorthand notation $s_\theta = \sin \theta$, $c_\theta = \cos \theta$, with θ defined as in eq. (2.17). The dots stand for potential gluon and W^\pm terms, for which there are no residual \mathcal{L}_6 corrections.

Eq. (2.18) shows that the contributions from Q_{HW} , Q_{HB} , Q_{HG} are fully reabsorbed in the definition of the fields and gauge couplings. As a consequence, these operators have no physical impact in the pure gauge sector, and they only contribute to Higgs-gauge interactions [63]. On the other hand, the operator Q_{HWB} introduces net modifications of all γ and Z couplings. In the former case the correction is a universal rescaling of the

electromagnetic constant, while in the latter case the corrections depend on the field's charges. In particular, in the Higgs case ($T_3 = -1/2, Q = 0$) this implies a correction $\propto \bar{C}_{HWB}$ to the Z mass term. The physical interpretation of these contributions requires defining a set of input observables and is deferred to section 4.

In `SMEFTsim`, the redefinitions described in this subsection are applied simultaneously at the Lagrangian level in the `FeynRules` model. The coupling constants' rescaling in eq. (2.11) is implemented in the replacement list `redefConst`. The field redefinitions are operated in the mass and charge eigenstate basis: the replacement list `rotateGaugeB` implements the net mismatch between the series of rotations (2.10), (2.13), (2.15) and the usual SM rotations, i.e.

$$G_\mu^a \rightarrow (1 + \bar{C}_{HG}) G_\mu^a, \quad (2.20)$$

$$\begin{pmatrix} W_\mu^+ \\ W_\mu^- \end{pmatrix} \rightarrow (1 + \bar{C}_{HW}) \begin{pmatrix} W_\mu^+ \\ W_\mu^- \end{pmatrix}, \quad (2.21)$$

$$\begin{pmatrix} Z_\mu \\ A_\mu \end{pmatrix} \rightarrow R_{ZA} \begin{pmatrix} Z_\mu \\ A_\mu \end{pmatrix}, \quad (2.22)$$

with

$$R_{ZA} = \begin{pmatrix} 1 + c_\theta^2 \bar{C}_{HW} + s_\theta^2 \bar{C}_{HB} + \frac{s_{2\theta}}{2} \bar{C}_{HWB} & \frac{s_{2\theta}}{2} (\bar{C}_{HW} - \bar{C}_{HB}) - \frac{c_{2\theta}}{2} \bar{C}_{HWB} + \frac{\Delta s_\theta^2}{s_{2\theta}} \\ \frac{s_{2\theta}}{2} (\bar{C}_{HW} - \bar{C}_{HB}) - \frac{c_{2\theta}}{2} \bar{C}_{HWB} - \frac{\Delta s_\theta^2}{s_{2\theta}} & 1 + c_\theta^2 \bar{C}_{HB} + s_\theta^2 \bar{C}_{HW} - \frac{s_{2\theta}}{2} \bar{C}_{HWB} \end{pmatrix}, \quad (2.23)$$

where

$$\Delta s_\theta^2 = \bar{C}_{HWB} g_1 g_W \frac{g_W^2 - g_1^2}{g_W^2 + g_1^2} = \frac{s_{4\theta}}{4} \bar{C}_{HWB} + \mathcal{O}(\Lambda^4), \quad (2.24)$$

is the correction to the mixing angle stemming from eq. (2.17).

3 Flavor assumptions

The SMEFT Lagrangian defined in section 1 is not invariant under flavor rotations of the fermion fields, so the latter should always be defined in order to avoid ambiguities. In `SMEFTsim`, the fields q, l, u, d, e are defined in the mass basis of the charged leptons and of the up-type quarks, in which the Yukawa matrices in eq. (1.5) take the form

$$Y_d \equiv Y_d^{(d)} V^\dagger, \quad Y_u \equiv Y_u^{(d)}, \quad Y_l \equiv Y_l^{(d)}. \quad (3.1)$$

The superscript (d) denotes diagonal matrices and V is the CKM matrix. This basis choice is consistently employed in the definition of both \mathcal{L}_{SM} and \mathcal{L}_6 , and for all the flavor assumptions implemented in `SMEFTsim`. The only special case are the `top` and `topU31` models, where quark mixing is entirely neglected by setting $V \equiv \mathbb{1}$.

Upon EWSB, the Lagrangian can be written in terms of the fermionic mass eigenstates. By definition the relation between the $SU(2)_L$ and mass bases is trivial for all fermion fields, except the left-handed quark doublet:⁶

$$q_p = \begin{pmatrix} u_{L,p} \\ V_{pr} d_{L,r} \end{pmatrix}, \quad l_p = \begin{pmatrix} \nu_{L,p} \\ e_{L,p} \end{pmatrix}, \quad u_p = u_{R,p}, \quad d_p = d_{R,p}, \quad e_p = e_{R,p}. \quad (3.2)$$

In unitary gauge, the relevant terms in the SM Lagrangian are therefore

$$\mathcal{L}_{\text{Fermions}} = -\frac{gW}{\sqrt{2}} W_\mu^+ \bar{u}_L \gamma^\mu V d_L - \frac{gW}{\sqrt{2}} W_\mu^+ \bar{\nu}_L \gamma^\mu e_L + \dots \quad (3.3)$$

$$\mathcal{L}_{\text{Yukawa}} = -\frac{v_T + h}{\sqrt{2}} \left[\bar{d}_R Y_d^{(d)} d_L + \bar{u}_R Y_u^{(d)} u_L + \bar{e}_R Y_l^{(d)} e_L \right] + \text{h.c.} \quad (3.4)$$

The CKM matrix is implemented in the Wolfenstein parameterization [64]:

$$V = \begin{pmatrix} 1 - \lambda_{CKM}^2/2 & \lambda_{CKM} & A \lambda_{CKM}^3 (\rho - i\eta) \\ -\lambda_{CKM} & 1 - \lambda_{CKM}^2/2 & A \lambda_{CKM}^2 \\ A \lambda_{CKM}^3 (1 - \rho - i\eta) & -A \lambda_{CKM}^2 & 1 \end{pmatrix}. \quad (3.5)$$

The numerical values employed for the parameters are listed in table 16.

SMEFTsim implements five alternative flavor scenarios: one with fully arbitrary indices, and four based on the implementation of different global symmetries. Three of these scenarios have been present since the first release, and two have been newly introduced in version 3.0. The following sub-sections review in detail the properties of the \mathcal{L}_6 operators within each setup and provide the corresponding parameter counting. A dictionary between the different flavor assumptions is provided in appendix C.

3.1 general: general flavor structure

Without further assumptions on the flavor structure of the SMEFT, \mathcal{L}_6 contains the operators in table 1, summed over all possible flavor combinations:

$$\begin{aligned} \mathcal{L}_6^{(5,6)} &= \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r=1}^3 C_{\alpha,pr} Q_{\alpha,pr} + \text{h.c.}, & \mathcal{L}_6^{(7)} &= \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r=1}^3 C_{\alpha,pr} Q_{\alpha,pr}, & (3.6) \\ \mathcal{L}_6^{(8a,8b,8c)} &= \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r,s,t=1}^3 C_{\alpha,prst} Q_{\alpha,prst}, & \mathcal{L}_6^{(8d)} &= \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r,s,t=1}^3 C_{\alpha,prst} Q_{\alpha,prst} + \text{h.c.} \end{aligned} \quad (3.7)$$

Not all flavor combinations included in this way are independent, due to intrinsic symmetry properties of the effective operators. SMEFTsim does not remove redundant terms from the sums in eqs. (3.6), (3.7). Instead, the symmetry relations are enforced in the definition of the tensor Wilson coefficients $C_{\alpha,pr(st)}$: only a minimum number of independent parameters is defined for each operator, as reported in appendix D, tables 18, 19, and all the entries of $C_{\alpha,pr(st)}$ are functions of these parameters, consistent with the relations described below.

⁶For economy of notation, we use the same letters u, d, e for the right-handed fields and for the mass eigenstates, both of them carrying flavor indices. To avoid ambiguities, the latter always carry L, R subscripts, while the former don't.

Classes 5 and 6. The operators in $\mathcal{L}_6^{(5,6)}$ are not Hermitian. Therefore each Wilson coefficient has 9 independent complex entries. In total, this gives 198 real parameters (counting independently real and imaginary parts).

Class 7. All operators in $\mathcal{L}_6^{(7)}$, except Q_{Hud} , are Hermitian. In this case, the diagonal entries of the Wilson coefficients are real, and the off-diagonal ones are related by

$$C_{pr} = C_{rp}^* . \tag{3.8}$$

In total, this class depends on 81 real parameters.

Class 8 a. All operators in $\mathcal{L}_6^{(8a)}$ are Hermitian. Moreover, each of the two currents that compose them is itself Hermitian. Therefore the following relation holds:

$$C_{prst} = C_{rpts} = C_{rpst}^* = C_{prts}^* . \tag{3.9}$$

In the operators $Q_{ll}, Q_{qq}^{(1)}, Q_{qq}^{(3)}$, the two currents contain the same fields, which leads to an additional exchange symmetry

$$C_{prst} = C_{stpr} . \tag{3.10}$$

Each of these three operators has then 15 real entries ($C_{pppp}, C_{pprr}, C_{prrp}$) and only 9 are independent, and 66 complex entries, 18 independent. Operators $Q_{lq}^{(1)}, Q_{lq}^{(3)}$ have each 9 real entries, all independent, and 72 complex ones, only 36 independent.

In total, this class depends on 297 real parameters.

Class 8 b. All operators in $\mathcal{L}_6^{(8b)}$ are Hermitian and composed of two Hermitian currents, so relation (3.9) holds for all Wilson coefficients in this class. Eq. (3.10) is valid in addition for C_{uu}, C_{dd}, C_{ee} .

The operator Q_{ee} is peculiar: because the e field is a singlet under both $SU(2)_L$ and $SU(3)_c$, this term is invariant under Fierz rearranging. This leads to the additional constraint

$$C_{prst} = C_{ptsr} . \tag{3.11}$$

The coefficient C_{ee} has then 15 real entries, 6 independent, and 66 complex entries, only 15 independent. The counting for the other operators is the same as for the invariants in class 8a, so $\mathcal{L}_6^{(8b)}$ has a total of 450 real parameters.

Class 8 c. All operators in $\mathcal{L}_6^{(8c)}$ are Hermitian and composed of two Hermitian currents, but no other symmetry is present. Therefore only relation (3.9) holds for all Wilson coefficients, leaving a total of 648 parameters.

Class 8 d. Finally, all operators in \mathcal{L}_6^{8d} are non-Hermitian. No symmetry relation is present and this class has 810 independent real parameters.

3.2 U35: maximal U(3)⁵ symmetry

The number of independent parameters is considerably reduced if a flavor symmetry is assumed. The maximal symmetry available for the SM fermion fields is the symmetry of the kinetic terms [65]: $U(3)^5 = U(3)_q \times U(3)_u \times U(3)_d \times U(3)_l \times U(3)_e$. Each field is assigned to a **3** representation of the associated group: denoting a generic $U(3)_\psi$ transformation by Ω_ψ , the transformation rules are [48]

$$q \mapsto \Omega_q q, \quad u \mapsto \Omega_u u, \quad d \mapsto \Omega_d d, \quad l \mapsto \Omega_l l, \quad e \mapsto \Omega_e e. \quad (3.12)$$

Vector currents $\bar{\psi}_p \gamma^\mu \psi_r$ are trivially made invariant under $U(3)^5$ by imposing a δ_{pr} contraction, that corresponds to the singlet composition of a $\bar{\mathbf{3}}$ and $\mathbf{3}$ representations. This is immediate to see applying the field transformations and using $\Omega_\psi \Omega_\psi^\dagger = \mathbb{1} = \Omega_\psi^\dagger \Omega_\psi$:

$$\bar{\psi} \gamma^\mu \psi \mapsto \bar{\psi} \Omega_\psi^\dagger \gamma^\mu \Omega_\psi \psi = \bar{\psi} \gamma^\mu \psi, \quad \psi = \{q, u, d, l, e\}. \quad (3.13)$$

Chirality-flipping currents, with either scalar or tensor Lorentz structure, violate the flavor symmetry. To permit the introduction of fermion masses, it is customary to promote the Yukawa couplings to *spurions* of the flavor symmetry, by assigning them transformation properties

$$Y_d \mapsto \Omega_d Y_d \Omega_q^\dagger, \quad Y_u \mapsto \Omega_u Y_u \Omega_q^\dagger, \quad Y_l \mapsto \Omega_e Y_l \Omega_l^\dagger. \quad (3.14)$$

In this way the structures

$$\bar{d} Y_d q^i, \quad \bar{u} Y_u q^i, \quad \bar{e} Y_l l^i, \quad (3.15)$$

are formally invariant. When the $U(3)^5$ symmetry is imposed, the flavor structure of each operator can be factored out of the Wilson coefficient, that becomes a scalar quantity:

$$\mathcal{L}_6^{(5,6)} = \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r=1}^3 C_\alpha X_{\alpha,pr} Q_{\alpha,pr} + \text{h.c.}, \quad \mathcal{L}_6^{(7)} = \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r=1}^3 C_\alpha X_{\alpha,pr} Q_{\alpha,pr}, \quad (3.16)$$

$$\mathcal{L}_6^{(8a,8b,8c)} = \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r,s,t=1}^3 C_\alpha X_{\alpha,prst} Q_{\alpha,prst}, \quad \mathcal{L}_6^{(8d)} = \frac{1}{\Lambda^2} \sum_\alpha \sum_{p,r,s,t=1}^3 C_\alpha X_{\alpha,prst} Q_{\alpha,prst} + \text{h.c.} \quad (3.17)$$

In the construction of the $U(3)^5$ symmetric Lagrangian, we do not define a power counting for insertions of the Yukawa couplings. Instead, we simply choose to retain the leading invariant structures for each operator, corresponding to no Yukawa insertions in $\mathcal{L}_6^{(7,8a,8b,8c)}$, one insertion in $\mathcal{L}_6^{(5,6)}$ and two insertions in C_{Hud} and $\mathcal{L}_6^{(8d)}$.

Classes 5 and 6. All the operators in $\mathcal{L}_6^{(5,6)}$ require the insertion of a Yukawa coupling:

$$\begin{aligned} X_{eH} &= X_{eW} = X_{eB} = Y_l^\dagger = Y_l^{(d)}, \\ X_{uH} &= X_{uW} = X_{uB} = X_{uG} = Y_u^\dagger = Y_u^{(d)}, \\ X_{dH} &= X_{dW} = X_{dB} = X_{dG} = Y_d^\dagger = V Y_d^{(d)}, \end{aligned} \quad (3.18)$$

where the last equality in each line holds in the up-quarks mass basis, eq. (3.1). Note that no net mixing among down-type quarks is induced in the mass basis, as the V in the spurion cancels against V^\dagger in the \bar{q} field, eq. (3.2). In fact, by construction all the operators in $\mathcal{L}_6^{(5,6)}$ have the same flavor structure as the SM Yukawas.

Because the operators are non-Hermitian, the associated C_α are complex. These classes therefore introduce 22 independent real parameters.

Class 7. All the currents appearing in the operators of class 7, except Q_{Hud} , are invariant under $U(3)$ with $X_\alpha = \mathbb{1}$. This implies that the flavor structure of this class is exactly the same as in the SM kinetic terms. For instance, the charged quark current induced by the operator $Q_{Hq}^{(3)}$ is aligned with the SM one (eq. (3.3)), that contains CKM mixing.⁷

In order to make Q_{Hud} invariant, it is necessary to insert the spurion product

$$X_{Hud} = Y_u Y_d^\dagger = Y_u^{(d)} V Y_d^{(d)}. \quad (3.19)$$

The number of independent real parameters is 9, as 7 out of 8 operators are Hermitian.

Class 8 a. Containing only vector currents, all the operators in $\mathcal{L}_6^{(8a)}$ are $U(3)^5$ invariant with the trivial flavor contraction $X_\alpha = \delta_{pr}\delta_{st}$.

The operators $Q_{ll}, Q_{qq}^{(1)}, Q_{qq}^{(3)}$ additionally admit the “crossed” contraction $X'_\alpha = \delta_{pt}\delta_{sr}$. This is an independent structure that cannot be arbitrarily rearranged into X_α : applying Fierz transformations in this case would introduce additional operators with $SU(2)_L$ triplet and $SU(3)_c$ octet contractions, see section 3.4. Therefore these operators are split into two invariants each, weighted by independent Wilson coefficients:

$$\Lambda^2 \mathcal{L}_6^{(8a)} = \left(C_{ll} \delta_{pr}\delta_{st} + C'_{ll} \delta_{pt}\delta_{sr} \right) Q_{ll,prst} + \left(C_{qq}^{(1)} \delta_{pr}\delta_{st} + C'_{qq} \delta_{pt}\delta_{sr} \right) Q_{qq,prst} + \dots \quad (3.20)$$

where the dots stand for contributions from the other operators in $\mathcal{L}_6^{(8a)}$.

As all operators are Hermitian, $\mathcal{L}_6^{(8a)}$ contains 8 real parameters.

Class 8 b. All operators in $\mathcal{L}_6^{(8b)}$ are invariant with $X_\alpha = \delta_{pr}\delta_{st}$.

The operators Q_{uu}, Q_{dd} additionally admit independent crossed contractions $X'_\alpha = \delta_{pt}\delta_{sr}$, and are treated analogously to $Q_{ll}, Q_{qq}^{(1),(3)}$. This is not the case for Q_{ee} that, as mentioned above, is invariant under Fierz rearrangements: in this particular case the two flavor contractions are equivalent. In total, there are 9 real parameters in $\mathcal{L}_6^{(8b)}$.

Class 8 c. All the operators in $\mathcal{L}_6^{(8c)}$ admit the invariant contraction is $X_\alpha = \delta_{pr}\delta_{st}$, leading to 8 independent real parameters.

Class 8 d. Finally, operators in $\mathcal{L}_6^{(8d)}$ require one Yukawa coupling insertion for each current. As they are not invariant under Fierz transformations, the operators $Q_{quqd}^{(1),(8)}$

⁷Note that this implies that, even though all Wilson coefficients are real, SM-sourced CP violation, due to the CKM phase in charged left-handed currents, is generally present in $\mathcal{L}_6^{(7)}$.

admit two independent contractions, mapped to one another by interchanging the two \bar{q} fields.⁸

$$\begin{aligned}
 X_{ledq} &= (Y_l^\dagger)_{pr} (Y_d)_{st} = (Y_l^{(d)})_{pr} (Y_d^{(d)} V^\dagger)_{st}, \\
 X_{quqd}^{(1)} &= X_{quqd}^{(8)} = (Y_u^\dagger)_{pr} (Y_d^\dagger)_{st} = (Y_u^{(d)})_{pr} (V Y_d^{(d)})_{st}, \\
 X_{quqd}^{(1)'} &= X_{quqd}^{(8)'} = (Y_u^\dagger)_{sr} (Y_d^\dagger)_{pt} = (Y_u^{(d)})_{sr} (V Y_d^{(d)})_{pt}, \\
 X_{lequ}^{(1)} &= X_{lequ}^{(3)} = (Y_l^\dagger)_{pr} (Y_u^\dagger)_{st} = (Y_l^{(d)})_{pr} (Y_u^{(d)})_{st}.
 \end{aligned} \tag{3.21}$$

Because the operators are non-Hermitian, there are 14 real parameters in $\mathcal{L}_6^{(8d)}$.

3.3 MFV: linear minimal flavor violation

The Minimal Flavor Violation ansatz [65, 67, 68] assumes that the only sources of flavor and CP violation in $\mathcal{L}_{\text{SMEFT}}$ are those already present in the SM, namely the Yukawa couplings and the CKM phase.

The requirement on CP violation implies that the Wilson coefficients of CP-odd bosonic operators scale with the Jarlskog invariant J [69, 70]:

$$\{C_{\tilde{G}}, C_{\tilde{W}}, C_{H\tilde{W}}, C_{H\tilde{B}}, C_{H\tilde{W}B}, C_{H\tilde{G}}\} \propto J, \tag{3.22}$$

$$J = \text{Im} \left(V_{pr} V_{st} V_{pt}^* V_{sr}^* \right) \simeq \eta A^2 \lambda_{CKM}^6 \left(1 - \frac{\lambda_{CKM}^2}{2} \right) \simeq 3 \times 10^{-5}. \tag{3.23}$$

As the J suppression is stronger, for instance, than a loop factor, these coefficients can be safely neglected within the scope of **SMEFTsim**. The corresponding operators are therefore not implemented in the MFV version. An analogous argument applies to sources of explicit CP violation in the fermion sector. In the Warsaw basis, these are the imaginary parts of the Wilson coefficients in \mathcal{L}_6 , that are not defined either in the **SMEFTsim** MFV models.

The requirement on flavor violation is realized imposing a $U(3)^5$ symmetry on the fermion fields and allowing for arbitrary $U(3)^5$ -invariant spurion insertions in the currents, that generate flavor violating effects. Such insertions are organized in an expansion in powers of the Yukawa couplings, that can be either resummed (obtaining a non-linear MFV formulation [71]) or treated as a truncated series. **SMEFTsim** adopts the latter option and retains contributions up to one power of Y_l and up to 3 powers of Y_u, Y_d .

The relevant spurion structures at this order are

$$S^u = Y_u Y_u^\dagger \sim (\mathbf{1}, \mathbf{8}, \mathbf{1}), \quad S^u \mapsto \Omega_u S^u \Omega_u^\dagger, \tag{3.24}$$

$$S^d = Y_d Y_d^\dagger \sim (\mathbf{1}, \mathbf{1}, \mathbf{8}), \quad S^d \mapsto \Omega_d S^d \Omega_d^\dagger, \tag{3.25}$$

$$S^{qu} = Y_u^\dagger Y_u \sim (\mathbf{8}, \mathbf{1}, \mathbf{1}), \quad S^{qu} \mapsto \Omega_q S^{qu} \Omega_q^\dagger, \tag{3.26}$$

$$S^{qd} = Y_d^\dagger Y_d \sim (\mathbf{8}, \mathbf{1}, \mathbf{1}), \quad S^{qd} \mapsto \Omega_q S^{qd} \Omega_q^\dagger. \tag{3.27}$$

The first column indicates the spurions' representation under the $U(3)_q \times U(3)_u \times U(3)_d$ group, while the second provides the corresponding transformation rules. All of them are

⁸The two X'_α structures (4 real parameters) for $Q_{quqd}^{(1),(8)}$ were not included in previous versions of **SMEFTsim**. I thank the authors of ref. [66] for pointing this out.

Hermitian and they satisfy

$$Y_u^\dagger S^u = S^{qu} Y_u^\dagger, \quad Y_u S^{qu} = S^u Y_u, \quad (3.28)$$

$$Y_d^\dagger S^d = S^{qd} Y_d^\dagger, \quad Y_d S^{qd} = S^d Y_d. \quad (3.29)$$

In the mass basis of the up quarks (eq. (3.1)) the spurions take the form

$$S^u \equiv (Y_u^{(d)})^2, \quad S^{qu} \equiv (Y_u^{(d)})^2 = S^u, \quad (3.30)$$

$$S^d \equiv (Y_d^{(d)})^2, \quad S^{qd} \equiv V(Y_d^{(d)})^2 V^\dagger. \quad (3.31)$$

Additional relevant structures in this basis are

$$S^{qu} Y_u^\dagger = (Y_u^{(d)})^3, \quad S^{qu} Y_d^\dagger = (Y_u^{(d)})^2 V Y_d^{(d)}, \quad (3.32)$$

$$S^{qd} Y_u^\dagger = V(Y_d^{(d)})^2 V^\dagger Y_u^{(d)}, \quad S^{qd} Y_d^\dagger = V(Y_d^{(d)})^3. \quad (3.33)$$

Classes 5 and 6. With the power counting chosen, $\mathcal{L}_6^{(5,6)}$ take the form

$$\begin{aligned} \Lambda^2 \mathcal{L}_6^{(5)} = & \sum_{p,r=1}^3 C_{eH} (Y_l^\dagger)_{pr} Q_{eH,pr} \\ & + \left[C_{uH}^{(0)} Y_u^\dagger + (\Delta^u C_{uH}) S^{qu} Y_u^\dagger + (\Delta^d C_{uH}) S^{qd} Y_u^\dagger \right]_{pr} Q_{uH,pr} \\ & + \left[C_{dH}^{(0)} Y_d^\dagger + (\Delta^u C_{dH}) S^{qu} Y_d^\dagger + (\Delta^d C_{dH}) S^{qd} Y_d^\dagger \right]_{pr} Q_{dH,pr} \\ & + \text{h.c.}, \end{aligned} \quad (3.34)$$

$$\begin{aligned} \Lambda^2 \mathcal{L}_6^{(6)} = & \sum_{p,r=1}^3 C_{eW} (Y_l^\dagger)_{pr} Q_{eW,pr} + C_{eB} (Y_l^\dagger)_{pr} Q_{eB,pr} \\ & + \left[C_{uG}^{(0)} Y_u^\dagger + (\Delta^u C_{uG}) S^{qu} Y_u^\dagger + (\Delta^d C_{uG}) S^{qd} Y_u^\dagger \right]_{pr} Q_{uG,pr} \\ & + \left[C_{uW}^{(0)} Y_u^\dagger + (\Delta^u C_{uW}) S^{qu} Y_u^\dagger + (\Delta^d C_{uW}) S^{qd} Y_u^\dagger \right]_{pr} Q_{uW,pr} \\ & + \left[C_{uB}^{(0)} Y_u^\dagger + (\Delta^u C_{uB}) S^{qu} Y_u^\dagger + (\Delta^d C_{uB}) S^{qd} Y_u^\dagger \right]_{pr} Q_{uB,pr} \\ & + \left[C_{dG}^{(0)} Y_d^\dagger + (\Delta^u C_{dG}) S^{qu} Y_d^\dagger + (\Delta^d C_{dG}) S^{qd} Y_d^\dagger \right]_{pr} Q_{dG,pr} \\ & + \left[C_{dW}^{(0)} Y_d^\dagger + (\Delta^u C_{dW}) S^{qu} Y_d^\dagger + (\Delta^d C_{dW}) S^{qd} Y_d^\dagger \right]_{pr} Q_{dW,pr} \\ & + \left[C_{dB}^{(0)} Y_d^\dagger + (\Delta^u C_{dB}) S^{qu} Y_d^\dagger + (\Delta^d C_{dB}) S^{qd} Y_d^\dagger \right]_{pr} Q_{dB,pr} \\ & + \text{h.c.}, \end{aligned} \quad (3.35)$$

where the parameters $C_\alpha^{(0)}$, $(\Delta^u C_\alpha)$, $(\Delta^d C_\alpha)$ are real, scalar quantities. The structures $Y_u^\dagger S^u$, $Y_d^\dagger S^d$ are also allowed for operators Q_{uX} , Q_{dX} respectively, but they are not independent due to eqs. (3.28), (3.29). These two classes contain a total of 27 real parameters.

Class 7. For the operators in class 7 we have

$$\begin{aligned}
 \Lambda^2 \mathcal{L}_6^{(7)} = & \sum_{p,r=1}^3 C_{Hl}^{(1)} \delta_{pr} Q_{Hl,pr}^{(1)} + C_{Hl}^{(3)} \delta_{pr} Q_{Hl,pr}^{(3)} + C_{He} \delta_{pr} Q_{He,pr} \\
 & + \left[C_{Hq}^{(1)(0)} \mathbb{1} + (\Delta^u C_{Hq}^{(1)}) S^{qu} + (\Delta^d C_{Hq}^{(1)}) S^{qd} \right]_{pr} Q_{Hq,pr}^{(1)} \\
 & + \left[C_{Hq}^{(3)(0)} \mathbb{1} + (\Delta^u C_{Hq}^{(3)}) S^{qu} + (\Delta^d C_{Hq}^{(3)}) S^{qd} \right]_{pr} Q_{Hq,pr}^{(3)} \\
 & + \left[C_{Hu}^{(0)} \mathbb{1} + (\Delta C_{Hu}) S^u \right]_{pr} Q_{Hu,pr} \\
 & + \left[C_{Hd}^{(0)} \mathbb{1} + (\Delta C_{Hd}) S^d \right]_{pr} Q_{Hd,pr} \\
 & + \left[\left[C_{Hud}^0 Y_u Y_d^\dagger \right]_{pr} Q_{Hud,pr} + \text{h.c.} \right].
 \end{aligned} \tag{3.36}$$

The total number of independent parameters in this class is 14.

Class 8 a. The operators of class 8a are composed of the same currents as those of class 7. MFV corrections have therefore an analogous structure. With the power counting chosen, the independent contractions are

$$\begin{aligned}
 \Lambda^2 \mathcal{L}_6^{(8a)} = & \sum_{p,r,s,t=1}^3 \left[C_{ll} \delta_{pr} \delta_{st} + C'_{ll} \delta_{pt} \delta_{sr} \right] Q_{ll,prst} \\
 & + \left[C_{qq}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{qq}^{(1)}) S_{pr}^{qu} \delta_{st} + (\Delta^d C_{qq}^{(1)}) S_{pr}^{qd} \delta_{st} + \right. \\
 & \quad \left. C_{qq}^{(1) \prime (0)} \delta_{pt} \delta_{sr} + (\Delta^u C_{qq}^{(1) \prime}) S_{pt}^{qu} \delta_{sr} + (\Delta^d C_{qq}^{(1) \prime}) S_{pt}^{qd} \delta_{sr} \right] Q_{qq,prst}^{(1)} \\
 & + \left[C_{qq}^{(3)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{qq}^{(3)}) S_{pr}^{qu} \delta_{st} + (\Delta^d C_{qq}^{(3)}) S_{pr}^{qd} \delta_{st} + \right. \\
 & \quad \left. C_{qq}^{(3) \prime (0)} \delta_{pt} \delta_{sr} + (\Delta^u C_{qq}^{(3) \prime}) S_{pt}^{qu} \delta_{sr} + (\Delta^d C_{qq}^{(3) \prime}) S_{pt}^{qd} \delta_{sr} \right] Q_{qq,prst}^{(3)} \\
 & + \left[C_{lq}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{lq}^{(1)}) \delta_{pr} S_{st}^{qu} + (\Delta^d C_{lq}^{(1)}) \delta_{pr} S_{st}^{qd} \right] Q_{lq,prst}^{(1)} \\
 & + \left[C_{lq}^{(3)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{lq}^{(3)}) \delta_{pr} S_{st}^{qu} + (\Delta^d C_{lq}^{(3)}) \delta_{pr} S_{st}^{qd} \right] Q_{lq,prst}^{(3)}.
 \end{aligned}$$

In the case of operators $Q_{ll}, Q_{qq}^{(1),(3)}$ two possible flavor contractions are allowed, as discussed in section 3.2. In $Q_{qq}^{(1),(3)}$ spurion insertions in the (st) , (sr) currents are redundant by symmetry. Class 8c contains 20 independent parameters.

Class 8 b. The MFV Lagrangian in class 8b has the form

$$\begin{aligned}
 \Lambda^2 \mathcal{L}_6^{(8b)} = & \sum_{p,r,s,t=1}^3 C_{ee} \delta_{pr} \delta_{st} Q_{ee,prst} \\
 & + \left[C_{uu}^{(0)} \delta_{pr} \delta_{st} + (\Delta C_{uu}) S_{pr}^u \delta_{st} + C_{uu}'^{(0)} \delta_{pt} \delta_{sr} + (\Delta C_{uu}') S_{pt}^u \delta_{sr} \right] Q_{uu,prst} \\
 & + \left[C_{dd}^{(0)} \delta_{pr} \delta_{st} + (\Delta C_{dd}) S_{pr}^d \delta_{st} + C_{dd}'^{(0)} \delta_{pt} \delta_{sr} + (\Delta C_{dd}') S_{pt}^d \delta_{sr} \right] Q_{dd,prst} \\
 & + \left[C_{eu}^{(0)} \delta_{st} + (\Delta C_{eu}) S_{st}^u \right] \delta_{pr} C_{eu,prst} \\
 & + \left[C_{ed}^{(0)} \delta_{st} + (\Delta C_{ed}) S_{st}^d \right] \delta_{pr} C_{ed,prst} \\
 & + \left[C_{ud}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{ud}^{(1)}) S_{pr}^u \delta_{st} + (\Delta^d C_{ud}^{(1)}) \delta_{pr} S_{st}^d \right] C_{ud,prst}^{(1)} \\
 & + \left[C_{ud}^{(8)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{ud}^{(8)}) S_{pr}^u \delta_{st} + (\Delta^d C_{ud}^{(8)}) \delta_{pr} S_{st}^d \right] C_{ud,prst}^{(8)}.
 \end{aligned} \tag{3.37}$$

As in the $U(3)^5$ symmetric case, Q_{uu}, Q_{dd} admit two independent flavor contractions, and spurion insertions in only one of their currents is required, by symmetry. Class 8b therefore contains a total of 19 independent parameters.

Class 8 c. For class 8c we have a total of 28 independent parameters:

$$\begin{aligned}
 \Lambda^2 \mathcal{L}_6^{(8c)} = & \sum_{p,r,s,t=1}^3 C_{le} \delta_{pr} \delta_{st} Q_{le,prst} \\
 & + \left[C_{lu}^{(0)} \delta_{st} + (\Delta C_{lu}) S_{st}^u \right] \delta_{pr} Q_{lu,prst} \\
 & + \left[C_{ld}^{(0)} \delta_{st} + (\Delta C_{ld}) S_{st}^d \right] \delta_{pr} Q_{ld,prst} \\
 & + \left[C_{qe}^{(0)} \delta_{pr} + (\Delta^u C_{qe}) S_{pr}^{qu} + (\Delta^d C_{qe}) S_{pr}^{qd} \right] \delta_{st} Q_{qe,prst} \\
 & + \left[C_{qu}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta_1^u C_{qu}^{(1)}) S_{pr}^{qu} \delta_{st} + (\Delta_1^d C_{qu}^{(1)}) S_{pr}^{qd} \delta_{st} + (\Delta_2 C_{qu}^{(1)}) \delta_{pr} S_{st}^u + \right. \\
 & \quad \left. C_{qu}^{(1) \prime(0)} (Y_u^\dagger)_{pt} (Y_u)_{sr} \right] Q_{qu,prst}^{(1)} \\
 & + \left[C_{qu}^{(8)(0)} \delta_{pr} \delta_{st} + (\Delta_1^u C_{qu}^{(8)}) S_{pr}^{qu} \delta_{st} + (\Delta_1^d C_{qu}^{(8)}) S_{pr}^{qd} \delta_{st} + (\Delta_2 C_{qu}^{(8)}) \delta_{pr} S_{st}^u + \right. \\
 & \quad \left. C_{qu}^{(8) \prime(0)} (Y_u^\dagger)_{pt} (Y_u)_{sr} \right] Q_{qu,prst}^{(8)} \\
 & + \left[C_{qd}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta_1^u C_{qd}^{(1)}) S_{pr}^{qu} \delta_{st} + (\Delta_1^d C_{qd}^{(1)}) S_{pr}^{qd} \delta_{st} + (\Delta_2 C_{qd}^{(1)}) \delta_{pr} S_{st}^d + \right. \\
 & \quad \left. C_{qd}^{(1) \prime(0)} (Y_d^\dagger)_{pt} (Y_d)_{sr} \right] Q_{qd,prst}^{(1)} \\
 & + \left[C_{qd}^{(8)(0)} \delta_{pr} \delta_{st} + (\Delta_1^u C_{qd}^{(8)}) S_{pr}^{qu} \delta_{st} + (\Delta_1^d C_{qd}^{(8)}) S_{pr}^{qd} \delta_{st} + (\Delta_2 C_{qd}^{(8)}) \delta_{pr} S_{st}^d + \right. \\
 & \quad \left. C_{qd}^{(8) \prime(0)} (Y_d^\dagger)_{pt} (Y_d)_{sr} \right] Q_{qd,prst}^{(8)}.
 \end{aligned}$$

Note that the operators $O_{qu}^{(1),(8)}, Q_{qd}^{(1),(8)}$ admit a contraction $[pt][sr]$ with a Yukawa insertion in each current.

Class 8 d. Finally, the operators of class 8d have the structure

$$\begin{aligned}
\Lambda^2 \mathcal{L}_6^{(8d)} = & \sum_{p,r,s,t=1}^3 \left[C_{ledq}^{(0)} Y_{d,st} + (\Delta^u C_{ledq}) (Y_d S^{qu})_{st} + (\Delta^d C_{ledq}) (Y_d S^{qd})_{st} \right] (Y_l^\dagger)_{pr} Q_{ledq,prst} \\
& + \left[C_{quqd}^{(1)} Y_{u,pr}^\dagger Y_{d,st}^\dagger + C_{quqd}^{(1)'} Y_{u,sr}^\dagger Y_{d,pt}^\dagger \right] Q_{quqd,prst}^{(1)} \\
& + \left[C_{quqd}^{(8)} Y_{u,pr}^\dagger Y_{d,st}^\dagger + C_{quqd}^{(8)'} Y_{u,sr}^\dagger Y_{d,pt}^\dagger \right] Q_{quqd,prst}^{(8)} \\
& + \left[C_{lequ}^{(1)(0)} Y_{u,st}^\dagger + (\Delta^u C_{lequ}^{(1)}) (S^{qu} Y_u^\dagger)_{st} + (\Delta^d C_{lequ}^{(1)}) (S^{qd} Y_u^\dagger)_{st} \right] (Y_l^\dagger)_{pr} Q_{lequ,prst}^{(1)} \\
& + \left[C_{lequ}^{(3)(0)} Y_{u,st}^\dagger + (\Delta^u C_{lequ}^{(3)}) (S^{qu} Y_u^\dagger)_{st} + (\Delta^d C_{lequ}^{(3)}) (S^{qd} Y_u^\dagger)_{st} \right] (Y_l^\dagger)_{pr} Q_{lequ,prst}^{(3)} \\
& + \text{h.c.},
\end{aligned} \tag{3.38}$$

The total number of parameters is 13.

3.4 top, topU31: $U(2)^3$ symmetry in the quark sector

Two new sets of models have been introduced in version 3.0, that implement a flavor structure consistent with the recommendations of ref. [36] for the SMEFT interpretation of top quark measurements. The formalism builds upon [72–74] and is defined by the following assumptions:

- quarks of the first two generations and quarks of the 3rd are described by independent fields. We denote them respectively by (q_p, u_p, d_p) with $p = \{1, 2\}$ and by (Q, t, b) .
- a symmetry $U(2)^3 = U(2)_q \times U(2)_u \times U(2)_d$ is imposed on the Lagrangian, under which only the light quarks transform:

$$q \mapsto \Omega_q q, \quad u \mapsto \Omega_u u, \quad d \mapsto \Omega_d d, \quad Q \mapsto Q, \quad t \mapsto t, \quad b \mapsto b. \tag{3.39}$$

- mixing effects in the quark sector are neglected and $V_{CKM} \equiv \mathbb{1}$ is assumed.

This choice greatly simplifies the structure of the Lagrangian, as mixing between the light and heavy quarks can only be introduced through extra $U(2)$ spurions [66, 72].

With this notation, the SM Lagrangian is

$$\begin{aligned}
\mathcal{L}_{\text{fermions}} = & i\bar{q}\not{D}q + i\bar{u}\not{D}u + i\bar{d}\not{D}d + i\bar{Q}\not{D}Q + i\bar{t}\not{D}t + i\bar{b}\not{D}b \quad + \text{leptons}, \\
\mathcal{L}_{\text{Yukawa}} = & -\bar{d}Y_d H^\dagger q - \bar{u}Y_u \tilde{H}^\dagger q - y_b \bar{b} H^\dagger Q - y_t \bar{t} \tilde{H}^\dagger Q \quad + \text{leptons},
\end{aligned} \tag{3.40}$$

with the Yukawas of the light quarks $Y_u \equiv \text{diag}(y_u, y_c)$, $Y_d \equiv \text{diag}(y_d, y_s)$ promoted to spurions of $U(2)$

$$Y_u \mapsto \Omega_u Y_u \Omega_q^\dagger, \quad Y_d \mapsto \Omega_d Y_d \Omega_q^\dagger, \tag{3.41}$$

while y_t, y_b do not transform under any symmetry. As a consequence, only $(\bar{L}R), (\bar{R}L)$ currents with light quarks need to be weighted by Yukawa insertions.

It is convenient to construct a $U(2)^3$ invariant basis mapping the fermionic operators of table 1 to the notation with 6 quark fields. We choose the set given in table 2, where,

analogously to the $U(3)^5$ case, we retain the least Yukawa-suppressed $U(2)^3$ -invariant contractions for each operator in the Warsaw basis.

In the lepton sector we consider two alternative ansätze:

(a) a $U(1)_{l+e}^3 = U(1)_e \times U(1)_\mu \times U(1)_\tau$ symmetry under which the fields transform as

$$l_1 \mapsto e^{i\alpha_e} l_1, \quad l_2 \mapsto e^{i\alpha_\mu} l_2, \quad l_3 \mapsto e^{i\alpha_\tau} l_3, \quad (3.42)$$

$$e_1 \mapsto e^{i\alpha_e} e_1, \quad e_2 \mapsto e^{i\alpha_\mu} e_2, \quad e_3 \mapsto e^{i\alpha_\tau} e_3. \quad (3.43)$$

This matches the “baseline” scenario in ref. [36] and corresponds to simple flavor-diagonality. It is implemented in the **top** models.

(b) a $U(3)^2 = U(3)_l \times U(3)_e$ symmetry under which

$$l \mapsto \Omega_l l, \quad e \mapsto \Omega_e e, \quad Y_l \mapsto \Omega_e Y_l \Omega_l^\dagger. \quad (3.44)$$

In the lepton sector, this setup matches exactly the structure of the **U35** and **MFV** models. It is more restrictive compared to $U(1)_{l+e}^3$ and contains fewer free parameters.

It is implemented in the **topU31** models.

In the $U(1)_{l+e}^3$ symmetric case, no transformation rule needs to be assigned to Y_l , as left- and right-handed leptons transform under the same symmetry. This implies that $(\bar{L}R)$, $(\bar{R}L)$ lepton currents are weighted by Y_l in the **topU31** models but not in the **top** ones.

Classes 5 and 6. The basis of quark operators for $\mathcal{L}_6^{(5)}$ and $\mathcal{L}_6^{(6)}$ in table 2 is easily constructed splitting the quark currents for the first 2 and the 3rd generations. Insertions of Y_u^\dagger, Y_d^\dagger in light quark currents, that are required for $U(2)^3$ invariance, are embedded in the operator definitions. $\mathcal{L}_6^{(5,6)}$ contain in total 32 real parameters (16 complex) coming from quark invariants.

When $U(1)_{l+e}^3$ is imposed (**top** models) on the lepton fields, $Q_{eH,pr}, Q_{eW,pr}, Q_{eB,pr}$ admit 3 independent contractions each, one per generation. When the more restrictive $U(3)^2$ is imposed (**topU31** models), each operator is associated to only one complex Wilson coefficient.

The total number of real independent parameters in $\mathcal{L}_6^{(5,6)}$ is therefore 50 in the **top** case and 38 in the **topU31** case. The Lagrangian is

$$\begin{aligned} \Lambda^2 \mathcal{L}_6^{(5)} &= C_{uH} Q_{uH} + C_{tH} Q_{tH} + C_{dH} Q_{dH} + C_{bH} Q_{bH} \\ &+ \begin{cases} \sum_{p=1}^3 (C_{eH})_{pp} Q_{eH,pp} & U(1)_{l+e}^3 \text{ [top]} \\ \sum_{p,r=1}^3 C_{eH} (Y_l^\dagger)_{pr} Q_{eH,pr} & U(3)^2 \text{ [topU31]} \end{cases} \\ &+ \text{h.c.}, \end{aligned} \quad (3.45)$$

$$\begin{aligned} \Lambda^2 \mathcal{L}_6^{(6)} &= C_{uW} Q_{uW} + C_{uB} Q_{uB} + C_{uG} Q_{uG} + C_{tW} Q_{tW} + C_{tB} Q_{tB} + C_{tG} Q_{tG} \\ &+ C_{dW} Q_{dW} + C_{dB} Q_{dB} + C_{dG} Q_{dG} + C_{bW} Q_{bW} + C_{bB} Q_{bB} + C_{bG} Q_{bG} \\ &+ \begin{cases} \sum_{p,r=1}^3 (C_{eW})_{pp} Q_{eW,pp} + (C_{eB})_{pp} Q_{eB,pp} & U(1)_{l+e}^3 \text{ [top]} \\ \sum_{p,r=1}^3 C_{eW} (Y_l^\dagger)_{pr} Q_{eW,pr} + C_{eB} (Y_l^\dagger)_{pr} Q_{eB,pr} & U(3)^2 \text{ [topU31]} \end{cases} \\ &+ \text{h.c.} \end{aligned} \quad (3.46)$$

$\mathcal{L}_6^{(5)} - \psi^2 H^3$							
Q_{uH}	$(H^\dagger H)(\bar{q} Y_u^\dagger u \tilde{H})$	Q_{dH}	$(H^\dagger H)(\bar{q} Y_d^\dagger d H)$	Q_{eH}	$(H^\dagger H)(\bar{l}_p e_r H)$		
Q_{tH}	$(H^\dagger H)(\bar{Q} \tilde{H} t)$	Q_{bH}	$(H^\dagger H)(\bar{Q} H b)$				
$\mathcal{L}_6^{(6)} - \psi^2 XH$							
Q_{eW}	$(\bar{l}_p \sigma^{\mu\nu} e_r) \sigma^i H W_{\mu\nu}^i$	Q_{uW}	$(\bar{q} Y_u^\dagger \sigma^{\mu\nu} u) \sigma^i \tilde{H} W_{\mu\nu}^i$	Q_{uB}	$(\bar{q} Y_u^\dagger \sigma^{\mu\nu} u) \tilde{H} B_{\mu\nu}$	Q_{uG}	$(\bar{q} Y_u^\dagger \sigma^{\mu\nu} T^a u) \tilde{H} G_{\mu\nu}^a$
Q_{eB}	$(\bar{l}_p \sigma^{\mu\nu} e_r) H B_{\mu\nu}$	Q_{tW}	$(\bar{Q} \sigma^{\mu\nu} t) \sigma^i \tilde{H} W_{\mu\nu}^i$	Q_{tB}	$(\bar{Q} \sigma^{\mu\nu} t) \tilde{H} B_{\mu\nu}$	Q_{tG}	$(\bar{Q} \sigma^{\mu\nu} T^a t) \tilde{H} G_{\mu\nu}^a$
Q_{dW}	$(\bar{q} Y_d^\dagger \sigma^{\mu\nu} d) \sigma^i H W_{\mu\nu}^i$	Q_{dB}	$(\bar{q} Y_d^\dagger \sigma^{\mu\nu} d) H B_{\mu\nu}$	Q_{dG}	$(\bar{q} Y_d^\dagger \sigma^{\mu\nu} T^a d) H G_{\mu\nu}^a$		
Q_{bW}	$(\bar{Q} \sigma^{\mu\nu} b) \sigma^i H W_{\mu\nu}^i$	Q_{bB}	$(\bar{Q} \sigma^{\mu\nu} b) H B_{\mu\nu}$	Q_{bG}	$(\bar{Q} \sigma^{\mu\nu} T^a b) H G_{\mu\nu}^a$		
$\mathcal{L}_6^{(7)} - \psi^2 H^2 D$							
$Q_{Hl}^{(1)}$	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{l}_p \gamma^\mu l_r)$	$Q_{Hl}^{(3)}$	$(H^\dagger i \overleftrightarrow{D}_\mu^i H)(\bar{l}_p \sigma^i \gamma^\mu l_r)$	Q_{He}	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{e}_p \gamma^\mu e_r)$		
$Q_{Hq}^{(1)}$	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{q} \gamma^\mu q)$	$Q_{Hq}^{(3)}$	$(H^\dagger i \overleftrightarrow{D}_\mu^i H)(\bar{q} \sigma^i \gamma^\mu q)$	Q_{Hu}	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{u} \gamma^\mu u)$	Q_{Hd}	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{d} \gamma^\mu d)$
$Q_{HQ}^{(1)}$	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{Q} \gamma^\mu Q)$	$Q_{HQ}^{(3)}$	$(H^\dagger i \overleftrightarrow{D}_\mu^i H)(\bar{Q} \sigma^i \gamma^\mu Q)$	Q_{Ht}	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{t} \gamma^\mu t)$	Q_{Hb}	$(H^\dagger i \overleftrightarrow{D}_\mu H)(\bar{b} \gamma^\mu b)$
Q_{Hud}	$i(\tilde{H}^\dagger D_\mu H)(\bar{u} Y_u Y_d^\dagger \gamma^\mu d)$	Q_{Htb}	$i(\tilde{H}^\dagger D_\mu H)(\bar{t} \gamma^\mu b)$				
$\mathcal{L}_6^{(8a)} - (\bar{L}L)(\bar{L}L)$							
$Q_{lq}^{(1)}$	$(\bar{l}_p \gamma_\mu l_r)(\bar{q} \gamma^\mu q)$	$Q_{lq}^{(3)}$	$(\bar{l}_p \sigma^i \gamma_\mu l_r)(\bar{q} \sigma^i \gamma^\mu q)$	Q_{ll}	$(\bar{l}_p \gamma_\mu l_r)(\bar{l}_s \gamma^\mu l_t)$		
$Q_{lQ}^{(1)}$	$(\bar{l}_p \gamma_\mu l_r)(\bar{Q} \gamma^\mu Q)$	$Q_{lQ}^{(3)}$	$(\bar{l}_p \sigma^i \gamma_\mu l_r)(\bar{Q} \sigma^i \gamma^\mu Q)$	$Q_{lQ}^{(1)}$	$(\bar{Q} \gamma_\mu Q)(\bar{Q} \gamma^\mu Q)$	$Q_{lQ}^{(8)}$	$(\bar{Q} T^a \gamma_\mu Q)(\bar{Q} T^a \gamma^\mu Q)$
$Q_{qq}^{(1,1)}$	$(\bar{q} \gamma_\mu q)(\bar{q} \gamma^\mu q)$	$Q_{qq}^{(1,8)}$	$(\bar{q} T^a \gamma_\mu q)(\bar{q} T^a \gamma^\mu q)$	$Q_{qq}^{(3,1)}$	$(\bar{q} \sigma^i \gamma_\mu q)(\bar{q} \sigma^i \gamma^\mu q)$	$Q_{qq}^{(3,8)}$	$(\bar{q} \sigma^i T^a \gamma_\mu q)(\bar{q} \sigma^i T^a \gamma^\mu q)$
$Q_{Qq}^{(1,1)}$	$(\bar{Q} \gamma_\mu Q)(\bar{q} \gamma^\mu q)$	$Q_{Qq}^{(1,8)}$	$(\bar{Q} T^a \gamma_\mu Q)(\bar{q} T^a \gamma^\mu q)$	$Q_{Qq}^{(3,1)}$	$(\bar{Q} \sigma^i \gamma_\mu Q)(\bar{q} \sigma^i \gamma^\mu q)$	$Q_{Qq}^{(3,8)}$	$(\bar{Q} \sigma^i T^a \gamma_\mu Q)(\bar{q} \sigma^i T^a \gamma^\mu q)$
$\mathcal{L}_6^{(8b)} - (\bar{R}R)(\bar{R}R)$							
Q_{eu}	$(\bar{e}_p \gamma_\mu e_r)(\bar{u} \gamma^\mu u)$	Q_{ed}	$(\bar{e}_p \gamma_\mu e_r)(\bar{d} \gamma^\mu d)$	Q_{ee}	$(\bar{e}_p \gamma_\mu e_r)(\bar{e}_s \gamma^\mu e_t)$		
Q_{et}	$(\bar{e}_p \gamma_\mu e_r)(\bar{t} \gamma^\mu t)$	Q_{eb}	$(\bar{e}_p \gamma_\mu e_r)(\bar{b} \gamma^\mu b)$	Q_{tt}	$(\bar{t} \gamma_\mu t)(\bar{t} \gamma^\mu t)$	Q_{bb}	$(\bar{b} \gamma_\mu b)(\bar{b} \gamma^\mu b)$
$Q_{uu}^{(1)}$	$(\bar{u} \gamma_\mu u)(\bar{u} \gamma^\mu u)$	$Q_{uu}^{(8)}$	$(\bar{u} T^a \gamma_\mu u)(\bar{u} T^a \gamma^\mu u)$	$Q_{tu}^{(1)}$	$(\bar{t} \gamma_\mu t)(\bar{u} \gamma^\mu u)$	$Q_{tu}^{(8)}$	$(\bar{t} T^a \gamma_\mu t)(\bar{u} T^a \gamma^\mu u)$
$Q_{dd}^{(1)}$	$(\bar{d} \gamma_\mu d)(\bar{d} \gamma^\mu d)$	$Q_{dd}^{(8)}$	$(\bar{d} T^a \gamma_\mu d)(\bar{d} T^a \gamma^\mu d)$	$Q_{db}^{(1)}$	$(\bar{b} \gamma_\mu b)(\bar{d} \gamma^\mu d)$	$Q_{db}^{(8)}$	$(\bar{b} T^a \gamma_\mu b)(\bar{d} T^a \gamma^\mu d)$
$Q_{ud}^{(1)}$	$(\bar{u} \gamma_\mu u)(\bar{d} \gamma^\mu d)$	$Q_{ud}^{(8)}$	$(\bar{u} T^a \gamma_\mu u)(\bar{d} T^a \gamma^\mu d)$	$Q_{td}^{(1)}$	$(\bar{t} \gamma_\mu t)(\bar{d} \gamma^\mu d)$	$Q_{td}^{(8)}$	$(\bar{t} T^a \gamma_\mu t)(\bar{d} T^a \gamma^\mu d)$
$Q_{ub}^{(1)}$	$(\bar{u} \gamma_\mu u)(\bar{b} \gamma^\mu b)$	$Q_{ub}^{(8)}$	$(\bar{u} T^a \gamma_\mu u)(\bar{b} T^a \gamma^\mu b)$	$Q_{tb}^{(1)}$	$(\bar{t} \gamma_\mu t)(\bar{b} \gamma^\mu b)$	$Q_{tb}^{(8)}$	$(\bar{t} T^a \gamma_\mu t)(\bar{b} T^a \gamma^\mu b)$
$Q_{utbd}^{(1)}$	$(Y_u Y_d^\dagger)_{pr}(\bar{u}_p \gamma_\mu t_r)(\bar{b} \gamma^\mu d_r)$	$Q_{utbd}^{(8)}$	$(Y_u Y_d^\dagger)_{pr}(\bar{u}_p T^a \gamma_\mu t_r)(\bar{b} T^a \gamma^\mu d_r)$				
$\mathcal{L}_6^{(8c)} - (\bar{L}L)(\bar{R}R)$							
Q_{lu}	$(\bar{l}_p \gamma_\mu l_r)(\bar{u} \gamma^\mu u)$	Q_{ld}	$(\bar{l}_p \gamma_\mu l_r)(\bar{d} \gamma^\mu d)$	Q_{qe}	$(\bar{q} \gamma_\mu q)(\bar{e}_p \gamma^\mu e_r)$	Q_{le}	$(\bar{l}_p \gamma_\mu l_r)(\bar{e}_s \gamma^\mu e_t)$
Q_{lt}	$(\bar{l}_p \gamma_\mu l_r)(\bar{t} \gamma^\mu t)$	Q_{lb}	$(\bar{l}_p \gamma_\mu l_r)(\bar{b} \gamma^\mu b)$	Q_{Qe}	$(\bar{Q} \gamma_\mu Q)(\bar{e}_p \gamma^\mu e_r)$		
$Q_{qu}^{(1)}$	$(\bar{q} \gamma_\mu q)(\bar{u} \gamma^\mu u)$	$Q_{Qu}^{(1)}$	$(\bar{Q} \gamma_\mu Q)(\bar{u} \gamma^\mu u)$	$Q_{qt}^{(1)}$	$(\bar{q} \gamma_\mu q)(\bar{t} \gamma^\mu t)$	$Q_{Qt}^{(1)}$	$(\bar{Q} \gamma_\mu Q)(\bar{t} \gamma^\mu t)$
$Q_{qu}^{(8)}$	$(\bar{q} T^a \gamma_\mu q)(\bar{u} T^a \gamma^\mu u)$	$Q_{Qu}^{(8)}$	$(\bar{Q} T^a \gamma_\mu Q)(\bar{u} T^a \gamma^\mu u)$	$Q_{qt}^{(8)}$	$(\bar{q} T^a \gamma_\mu q)(\bar{t} T^a \gamma^\mu t)$	$Q_{Qt}^{(8)}$	$(\bar{Q} T^a \gamma_\mu Q)(\bar{t} T^a \gamma^\mu t)$
$Q_{qd}^{(1)}$	$(\bar{q} \gamma_\mu q)(\bar{d} \gamma^\mu d)$	$Q_{Qd}^{(1)}$	$(\bar{Q} \gamma_\mu Q)(\bar{d} \gamma^\mu d)$	$Q_{qb}^{(1)}$	$(\bar{q} \gamma_\mu q)(\bar{b} \gamma^\mu b)$	$Q_{Qb}^{(1)}$	$(\bar{Q} \gamma_\mu Q)(\bar{b} \gamma^\mu b)$
$Q_{qd}^{(8)}$	$(\bar{q} T^a \gamma_\mu q)(\bar{d} T^a \gamma^\mu d)$	$Q_{Qd}^{(8)}$	$(\bar{Q} T^a \gamma_\mu Q)(\bar{d} T^a \gamma^\mu d)$	$Q_{qb}^{(8)}$	$(\bar{q} T^a \gamma_\mu q)(\bar{b} T^a \gamma^\mu b)$	$Q_{Qb}^{(8)}$	$(\bar{Q} T^a \gamma_\mu Q)(\bar{b} T^a \gamma^\mu b)$
$Q_{qQu}^{(1)}$	$(Y_u^\dagger)_{pr}(\bar{q}_p \gamma_\mu Q)(\bar{t}_r \gamma^\mu u_r)$	$Q_{qQu}^{(8)}$	$(Y_u^\dagger)_{pr}(\bar{q}_p T^a \gamma_\mu Q)(\bar{t}_r T^a \gamma^\mu u_r)$	$Q_{qQbd}^{(1)}$	$(Y_d^\dagger)_{pr}(\bar{q}_p \gamma_\mu Q)(\bar{b}_r \gamma^\mu d_r)$	$Q_{qQbd}^{(8)}$	$(Y_d^\dagger)_{pr}(\bar{q}_p T^a \gamma_\mu Q)(\bar{b}_r T^a \gamma^\mu d_r)$
$\mathcal{L}_6^{(8d)} - (\bar{L}R)(\bar{R}L), (\bar{L}R)(\bar{L}R)$							
Q_{leqj}	$(\bar{l}_p^\dagger e_r)(\bar{d} Y_d q_j)$	Q_{lebQ}	$(\bar{l}_p^\dagger e_r)(\bar{b} Q_j)$	$Q_{leQt}^{(1)}$	$(\bar{l}_p^\dagger e_r) \varepsilon_{jk} (\bar{Q}^k t)$	$Q_{leQt}^{(3)}$	$(\bar{l}_p^\dagger \sigma_{\mu\nu} e_r) \varepsilon_{jk} (\bar{Q}^k \sigma^{\mu\nu} t)$
$Q_{lequ}^{(1)}$	$(\bar{l}_p^\dagger e_r) \varepsilon_{jk} (\bar{q}^k Y_u^\dagger u)$	$Q_{lequ}^{(3)}$	$(\bar{l}_p^\dagger \sigma_{\mu\nu} e_r) \varepsilon_{jk} (\bar{q}^k Y_u^\dagger \sigma^{\mu\nu} u)$	$Q_{QtQb}^{(1)}$	$(\bar{Q}^j t) \varepsilon_{jk} (\bar{Q}^k b)$	$Q_{QtQb}^{(8)}$	$(\bar{Q}^j T^a t) \varepsilon_{jk} (\bar{Q}^k T^a b)$
$Q_{quqd}^{(1)}$	$(\bar{q}^\dagger Y_u^\dagger u) \varepsilon_{jk} (\bar{q}^k Y_d^\dagger d)$	$Q_{quqd}^{(8)}$	$(\bar{q}^\dagger Y_u^\dagger T^a u) \varepsilon_{jk} (\bar{q}^k Y_d^\dagger T^a d)$	$Q_{quqd}^{(1)'} $	$(Y_u^\dagger)_{sr} (Y_d^\dagger)_{pt} (\bar{q}_p^\dagger u_r) \varepsilon_{jk} (\bar{q}_s^k d_t)$	$Q_{quqd}^{(8)'}$	$(Y_u^\dagger)_{sr} (Y_d^\dagger)_{pt} (\bar{q}_p^\dagger T^a u_r) \varepsilon_{jk} (\bar{q}_s^k T^a d_t)$
$Q_{Qtqd}^{(1)}$	$(\bar{Q}^j t) \varepsilon_{jk} (\bar{q}^k Y_d^\dagger d)$	$Q_{Qtqd}^{(8)}$	$(\bar{Q}^j T^a t) \varepsilon_{jk} (\bar{q}^k Y_d^\dagger T^a d)$	$Q_{quQb}^{(1)}$	$(\bar{q}^\dagger Y_u^\dagger u) \varepsilon_{jk} (\bar{Q}^k b)$	$Q_{quQb}^{(8)}$	$(\bar{q}^\dagger Y_u^\dagger T^a u) \varepsilon_{jk} (\bar{Q}^k T^a b)$
$Q_{Quqb}^{(1)}$	$(Y_u^\dagger)_{pr} (\bar{Q}^j T^a u_r) \varepsilon_{jk} (\bar{q}_p^k b)$	$Q_{Quqb}^{(8)}$	$(Y_u^\dagger)_{pr} (\bar{Q}^j T^a u_r) \varepsilon_{jk} (\bar{q}_p^k T^a b)$	$Q_{qtQd}^{(1)}$	$(Y_d^\dagger)_{pr} (\bar{q}_p^\dagger t) \varepsilon_{jk} (\bar{Q}^k d_r)$	$Q_{qtQd}^{(8)}$	$(Y_d^\dagger)_{pr} (\bar{q}_p^\dagger T^a t) \varepsilon_{jk} (\bar{Q}^k T^a d_r)$

Table 2. Basis of fermionic operators for the top and topU31 flavor assumptions. Here (q, u, d) , Y_u, Y_d denote quarks of the first 2 generations and their 2×2 Yukawa matrices. Quark fields of the 3rd generation are (Q, t, b) . Flavor indices p, r, s, t run over $\{1, 2\}$ for light quarks and $\{1, 2, 3\}$ for leptons. Whenever flavor indices are not specified, they are implicitly contracted within each current.

Class 7. Class 7 depends on 12 real parameters from quark operators, plus 9 (3) real parameters from lepton operators in the **top** (**topU31**) case. Q_{Hud} is defined with a $Y_u Y_d^\dagger$ insertion to preserve $U(2)^3$, while Q_{Htb} is independent of the Yukawas.

$$\begin{aligned} \Lambda^2 \mathcal{L}_6^{(7)} = & C_{Hq}^{(1)} Q_{Hq}^{(1)} + C_{Hq}^{(3)} Q_{Hq}^{(3)} + C_{Hu} Q_{Hu} + C_{Hd} Q_{Hd} + [C_{Hud} Q_{Hud} + \text{h.c.}] \\ & + C_{HQ}^{(1)} Q_{HQ}^{(1)} + C_{HQ}^{(3)} Q_{HQ}^{(3)} + C_{Ht} Q_{Ht} + C_{Hb} Q_{Hb} + [C_{Htb} Q_{Htb} + \text{h.c.}] \quad (3.47) \\ & + \begin{cases} \sum_{p=1}^3 (C_{Hl}^{(1)})_{pp} Q_{Hl,pp}^{(1)} + (C_{Hl}^{(3)})_{pp} Q_{Hl,pp}^{(3)} + (C_{He})_{pp} Q_{He,pp}, & U(1)_{l+e}^3 [\text{top}] \\ \sum_{p,r=1}^3 C_{Hl}^{(1)} \delta_{pr} Q_{Hl,pr}^{(1)} + C_{Hl}^{(3)} \delta_{pr} Q_{Hl,pr}^{(3)} + C_{He} \delta_{pr} Q_{He,pr}, & U(3)^2 [\text{topU31}] \end{cases} \end{aligned}$$

Class 8 a. Class 8a contains 2 operators with 4 quarks. Mapping them to the formalism with 6 quark fields, each of them admits 5 independent $U(2)^3$ invariant contractions, that can be written

$$Q_{qq}^{(1)} = \sum_{p,r,s,t=1}^2 \delta_{pr} \delta_{st} (\bar{q}_p \gamma_\mu q_r) (\bar{q}_s \gamma^\mu q_t), \quad Q_{Qq}^{(1)} = \sum_{p,r=1}^2 \delta_{pr} (\bar{q}_p \gamma_\mu q_r) (\bar{Q} \gamma^\mu Q), \quad (3.48)$$

$$Q_{qq}^{(1)'} = \sum_{p,r,s,t=1}^2 \delta_{pt} \delta_{sr} (\bar{q}_p \gamma_\mu q_r) (\bar{q}_s \gamma^\mu q_t), \quad Q_{Qq}^{(1)'} = \sum_{p,t=1}^2 \delta_{pt} (\bar{q}_p \gamma_\mu Q) (\bar{Q} \gamma^\mu q_t), \quad (3.49)$$

$$Q_{QQ}^{(1)} = (\bar{Q} \gamma_\mu Q) (\bar{Q} \gamma^\mu Q), \quad (3.50)$$

and analogously for $Q_{qq}^{(3)}$. In practice, for analyses involving top quark processes it is convenient to trade ‘‘crossed’’ flavor contractions, as well as $Q_{QQ}^{(3)}$, for operators with a color octet structure. This is motivated by top processes being largely dominated by QCD interactions in the SM. The rotation is done using Fierz rearrangements and the completeness relations for $SU(2)$ and $SU(3)$

$$\sigma_{jk}^i \sigma_{mn}^i = 2\delta_{jn} \delta_{mk} - \delta_{jk} \delta_{mn}, \quad (3.51)$$

$$T_{ab}^A T_{cd}^A = \frac{1}{2} \delta_{ad} \delta_{cb} - \frac{1}{6} \delta_{ab} \delta_{cd}. \quad (3.52)$$

Consistent with the recommendations in ref. [36], **SMEFTsim** implements the invariants in table 2, that are related to those in eqs. (3.48)–(3.50) and their $Q_{qq}^{(3)}$ counterparts as:

$$Q_{QQ}^{(1)} = Q_{QQ}^{(1)}, \quad Q_{QQ}^{(3)} = -\frac{1}{3} Q_{QQ}^{(1)} + 4Q_{QQ}^{(8)}, \quad (3.53)$$

$$Q_{qq}^{(1)} = Q_{qq}^{(1,1)}, \quad Q_{qq}^{(1)'} = \frac{1}{6} (Q_{qq}^{(1,1)} + Q_{qq}^{(3,1)}) + Q_{qq}^{(1,8)} + Q_{qq}^{(3,8)}, \quad (3.54)$$

$$Q_{qq}^{(3)} = Q_{qq}^{(3,1)}, \quad Q_{qq}^{(3)'} = \frac{1}{2} Q_{qq}^{(1,1)} - \frac{1}{6} Q_{qq}^{(3,1)} + 3Q_{qq}^{(1,8)} - Q_{qq}^{(3,8)}, \quad (3.55)$$

$$Q_{Qq}^{(1)} = Q_{Qq}^{(1,1)}, \quad Q_{Qq}^{(1)'} = \frac{1}{6} (Q_{Qq}^{(1,1)} + Q_{Qq}^{(3,1)}) + Q_{Qq}^{(1,8)} + Q_{Qq}^{(3,8)}, \quad (3.56)$$

$$Q_{Qq}^{(3)} = Q_{Qq}^{(3,1)}, \quad Q_{Qq}^{(3)'} = \frac{1}{2} Q_{Qq}^{(1,1)} - \frac{1}{6} Q_{Qq}^{(3,1)} + 3Q_{Qq}^{(1,8)} - Q_{Qq}^{(3,8)}. \quad (3.57)$$

Eqs. (3.53)–(3.57) can be written compactly as $\vec{Q}_{\text{Warsaw}} = R \vec{Q}_{\text{top}}$, with \vec{Q}_{Warsaw} , \vec{Q}_{top} the two “operator vectors” and R a rotation matrix. The relation among the Wilson coefficients is then derived equating the Lagrangian written in the two bases:

$$\begin{aligned} \mathcal{L} &= \vec{C}_{\text{Warsaw}} \cdot \vec{Q}_{\text{Warsaw}} = (\vec{C}_{\text{Warsaw}})^T R \vec{Q}_{\text{top}} \\ &= \vec{C}_{\text{top}} \cdot \vec{Q}_{\text{top}}, \end{aligned} \quad (3.58)$$

with \vec{C}_{Warsaw} , \vec{C}_{top} the coefficients vectors. The solution is

$$\vec{C}_{\text{top}} = R^T \vec{C}_{\text{Warsaw}}, \quad \vec{C}_{\text{Warsaw}} = (R^T)^{-1} \vec{C}_{\text{top}}. \quad (3.59)$$

Explicitly:

$$C_{QQ}^{(1)} = C_{QQ}^{(1)} - \frac{1}{3} C_{QQ}^{(3)}, \quad C_{QQ}^{(8)} = 4C_{QQ}^{(3)}, \quad (3.60)$$

$$C_{qq}^{(1,1)} = C_{qq}^{(1)} + \frac{1}{6} C_{qq}^{(1)'} + \frac{1}{2} C_{qq}^{(3)'}, \quad C_{qq}^{(3,1)} = C_{qq}^{(3)} + \frac{1}{6} C_{qq}^{(1)'} - \frac{1}{6} C_{qq}^{(3)'}, \quad (3.61)$$

$$C_{qq}^{(1,8)} = C_{qq}^{(1)'} + 3C_{qq}^{(3)'}, \quad C_{qq}^{(3,8)} = C_{qq}^{(1)'} - C_{qq}^{(3)'}, \quad (3.62)$$

$$C_{Qq}^{(1,1)} = C_{Qq}^{(1)} + \frac{1}{6} C_{Qq}^{(1)'} + \frac{1}{2} C_{Qq}^{(3)'}, \quad C_{Qq}^{(3,1)} = C_{Qq}^{(3)} + \frac{1}{6} C_{Qq}^{(1)'} - \frac{1}{6} C_{Qq}^{(3)'}, \quad (3.63)$$

$$C_{Qq}^{(1,8)} = C_{Qq}^{(1)'} + 3C_{Qq}^{(3)'}, \quad C_{Qq}^{(3,8)} = C_{Qq}^{(1)'} - C_{Qq}^{(3)'}, \quad (3.64)$$

and the inverse

$$C_{QQ}^{(1)} = C_{QQ}^{(1)} + \frac{1}{12} C_{QQ}^{(8)}, \quad C_{QQ}^{(3)} = \frac{1}{4} C_{QQ}^{(8)}, \quad (3.65)$$

$$C_{qq}^{(1)} = C_{qq}^{(1,1)} - \frac{1}{6} C_{qq}^{(1,8)}, \quad C_{qq}^{(1)'} = \frac{1}{4} C_{qq}^{(1,8)} + \frac{3}{4} C_{qq}^{(3,8)}, \quad (3.66)$$

$$C_{qq}^{(3)} = C_{qq}^{(3,1)} - \frac{1}{6} C_{qq}^{(3,8)}, \quad C_{qq}^{(3)'} = \frac{1}{4} C_{qq}^{(1,8)} - \frac{1}{4} C_{qq}^{(3,8)}, \quad (3.67)$$

$$C_{Qq}^{(1)} = C_{Qq}^{(1,1)} - \frac{1}{6} C_{Qq}^{(1,8)}, \quad C_{Qq}^{(1)'} = \frac{1}{4} C_{Qq}^{(1,8)} + \frac{3}{4} C_{Qq}^{(3,8)}, \quad (3.68)$$

$$C_{Qq}^{(3)} = C_{Qq}^{(3,1)} - \frac{1}{6} C_{Qq}^{(3,8)}, \quad C_{Qq}^{(3)'} = \frac{1}{4} C_{Qq}^{(1,8)} - \frac{1}{4} C_{Qq}^{(3,8)}. \quad (3.69)$$

The operator Q_{ll} admits 2 independent contractions in the $U(3)^2$ symmetric case (eq. (3.20)) and 9 in the $U(1)_{l+e}^3$ case. We choose them as the $(C_{ll})_{prst}$ entries with indices $prst$ in the set

$$P_{ll} = \{1111, 2222, 3333, 1122, 1133, 2233, 1221, 1331, 2332\}. \quad (3.70)$$

Operators $Q_{lq}^{(1),(3)}$, $Q_{lQ}^{(1),(3)}$ admit 3 (1) contractions each in the top (topU31) case.

The Lagrangian is therefore

$$\begin{aligned}
 \Lambda^2 \mathcal{L}_6^{(8a)} = & C_{qq}^{(1,1)} Q_{qq}^{(1,1)} + C_{qq}^{(1,8)} Q_{qq}^{(1,8)} + C_{qq}^{(3,1)} Q_{qq}^{(3,1)} + C_{qq}^{(3,8)} Q_{qq}^{(3,8)} + C_{QQ}^{(1)} Q_{QQ}^{(1)} \\
 & + C_{Qq}^{(1,1)} Q_{Qq}^{(1,1)} + C_{Qq}^{(1,8)} Q_{Qq}^{(1,8)} + C_{Qq}^{(3,1)} Q_{Qq}^{(3,1)} + C_{Qq}^{(3,8)} Q_{Qq}^{(3,8)} + C_{QQ}^{(8)} Q_{QQ}^{(8)} \\
 & + \left\{ \begin{array}{l} \sum_{p=1}^3 (C_{lq}^{(1)})_{pp} Q_{lq,pp}^{(1)} + (C_{lq}^{(3)})_{pp} Q_{lq,pp}^{(3)} + (C_{lQ}^{(1)})_{pp} Q_{lQ,pp}^{(1)} + (C_{lQ}^{(3)})_{pp} Q_{lQ,pp}^{(3)} \\ + \sum_{prst \in P_U} (C_U)_{prst} Q_{ll,prst}, \end{array} \right. \quad \text{U}(1)_{l+e}^3 \text{ [top]} \\
 & + \left\{ \begin{array}{l} \sum_{p,r=1}^3 C_{lq}^{(1)} \delta_{pr} Q_{lq,pr}^{(1)} + C_{lq}^{(3)} \delta_{pr} Q_{lq,pr}^{(3)} + C_{lQ}^{(1)} \delta_{pr} Q_{lQ,pr}^{(1)} + C_{lQ}^{(3)} \delta_{pr} Q_{lQ,pr}^{(3)} \\ + \sum_{p,r,s,t=1}^3 (C_U \delta_{pr} \delta_{st} + C'_U \delta_{pt} \delta_{sr}) Q_{ll,prst}, \end{array} \right. \quad \text{U}(3)^2 \text{ [topU31]}
 \end{aligned} \tag{3.71}$$

Note that the allowed flavor contractions in the $\text{U}(1)_{l+r}^3$ and $\text{U}(3)^2$ cases are the same, but the different symmetry properties generally lead to different relative normalizations. For instance, considering the (1111) and (1122) entries, one has

$$\Lambda^2 \mathcal{L}_6^{(8a)} \supset \left\{ \begin{array}{ll} (C_U)_{1111} Q_{1111} + (C_U)_{1122} Q_{1122} + \dots & \text{U}(1)_{l+e}^3 \text{ [top]} \\ (C_U + C'_U) Q_{1111} + 2C_U Q_{1122} + \dots & \text{U}(3)^2 \text{ [topU31]} \end{array} \right. \tag{3.72}$$

where the relative 2 between the C_U contributions to Q_{1122} and Q_{1111} is due to $\text{U}(3)^2$ requiring to sum over both the 1122 and 2211 contractions, that are equivalent for this particular operator. In total, $\mathcal{L}_6^{(8a)}$ contains 31 independent real parameters in the **top** case and 16 in the **topU31** case.

Class 8 b. A basis rotation analogous to the one performed in $\mathcal{L}_6^{(8a)}$ is applied to Q_{uu}, Q_{ud} in $\mathcal{L}_6^{(8b)}$. No modification is needed for $Q_{ud}^{(1),(8)}$ as in this case the color octet contraction is already manifest. The set of 5 independent $\text{U}(2)^3$ -invariant contractions in the Warsaw basis is in this case

$$Q_{uu} = \sum_{p,r,s,t=1}^2 \delta_{pr} \delta_{st} (\bar{u}_p \gamma_\mu u_r) (\bar{u}_s \gamma^\mu u_t), \quad Q_{tu} = \sum_{p,r=1}^2 \delta_{pr} (\bar{u}_p \gamma_\mu u_r) (\bar{t} \gamma^\mu t), \tag{3.73}$$

$$Q'_{uu} = \sum_{p,r,s,t=1}^2 \delta_{pt} \delta_{sr} (\bar{u}_p \gamma_\mu u_r) (\bar{u}_s \gamma^\mu u_t), \quad Q'_{tu} = \sum_{p,t=1}^2 \delta_{pt} (\bar{u}_p \gamma_\mu t) (\bar{t} \gamma^\mu u_t), \tag{3.74}$$

$$Q_{tt} = (\bar{t} \gamma_\mu t) (\bar{t} \gamma^\mu t), \tag{3.75}$$

and analogously for the Q_{dd} counterparts. Using Fierz transformations and eqs. (3.51), (3.52):

$$Q_{tt} = Q_{tt}, \quad Q_{bb} = Q_{bb}, \tag{3.76}$$

$$Q_{uu} = Q_{uu}^{(1)}, \quad Q'_{uu} = \frac{1}{3} Q_{uu}^{(1)} + 2Q_{uu}^{(8)}, \tag{3.77}$$

$$Q_{dd} = Q_{dd}^{(1)}, \quad Q'_{dd} = \frac{1}{3} Q_{dd}^{(1)} + 2Q_{dd}^{(8)}, \tag{3.78}$$

$$Q_{tu} = Q_{tu}^{(1)}, \quad Q'_{tu} = \frac{1}{3} Q_{tu}^{(1)} + 2Q_{tu}^{(8)}, \tag{3.79}$$

$$Q_{bd} = Q_{bd}^{(1)}, \quad Q'_{bd} = \frac{1}{3} Q_{bd}^{(1)} + 2Q_{bd}^{(8)}, \tag{3.80}$$

where the operators on the right-hand side of the equations are defined in table 2. The relations among Wilson coefficients are

$$C_{tt} = C_{tt}, \quad C_{bb} = C_{bb}, \quad (3.81)$$

$$C_{uu}^{(1)} = C_{uu} + \frac{1}{3}C'_{uu}, \quad C_{uu}^{(8)} = 2C'_{uu}, \quad (3.82)$$

$$C_{dd}^{(1)} = C_{dd} + \frac{1}{3}C'_{dd}, \quad C_{dd}^{(8)} = 2C'_{dd}, \quad (3.83)$$

$$C_{tu}^{(1)} = C_{tu} + \frac{1}{3}C'_{tu}, \quad C_{tu}^{(8)} = 2C'_{tu}, \quad (3.84)$$

$$C_{bd}^{(1)} = C_{bd} + \frac{1}{3}C'_{bd}, \quad C_{bd}^{(8)} = 2C'_{bd}, \quad (3.85)$$

and the inverse

$$C_{tt} = C_{tt}, \quad C_{bb} = C_{bb}, \quad (3.86)$$

$$C_{uu} = C_{uu}^{(1)} - \frac{1}{6}C_{uu}^{(8)}, \quad C'_{uu} = \frac{1}{2}C_{uu}^{(8)}, \quad (3.87)$$

$$C_{dd} = C_{dd}^{(1)} - \frac{1}{6}C_{dd}^{(8)}, \quad C'_{dd} = \frac{1}{2}C_{dd}^{(8)}, \quad (3.88)$$

$$C_{tu} = C_{tu}^{(1)} - \frac{1}{6}C_{tu}^{(8)}, \quad C'_{tu} = \frac{1}{2}C_{tu}^{(8)}, \quad (3.89)$$

$$C_{bd} = C_{bd}^{(1)} - \frac{1}{6}C_{bd}^{(8)}, \quad C'_{bd} = \frac{1}{2}C_{bd}^{(8)}. \quad (3.90)$$

The operator Q_{ee} admits 6 independent contractions in the **top** case, with indices that we choose in the set

$$P_{ee} = \{1111, 2222, 3333, 1122, 1133, 2233\}. \quad (3.91)$$

In the $U(3)^2$ case, there is instead only 1 available contraction. Each of the operators $Q_{lu}^{(1),(3)}$, $Q_{ld}^{(1),(3)}$ is mapped into 6 (2) independent invariants in the **top** (**topU31**) case. The Lagrangian for class 8a has the form

$$\begin{aligned} \Lambda^2 \mathcal{L}_6^{(8b)} = & C_{uu}^{(1)} Q_{uu}^{(1)} + C_{uu}^{(8)} Q_{uu}^{(8)} + C_{tt} Q_{tt} + C_{tu}^{(1)} Q_{tu}^{(1)} + C_{tu}^{(8)} Q_{tu}^{(8)} \\ & + C_{dd}^{(1)} Q_{dd}^{(1)} + C_{dd}^{(8)} Q_{dd}^{(8)} + C_{bb} Q_{bb} + C_{bd}^{(1)} Q_{bd}^{(1)} + C_{bd}^{(8)} Q_{bd}^{(8)} \\ & + C_{ud}^{(1)} Q_{ud}^{(1)} + C_{ud}^{(8)} Q_{ud}^{(8)} + C_{td}^{(1)} Q_{td}^{(1)} + C_{td}^{(8)} Q_{td}^{(8)} + C_{ub}^{(1)} Q_{ub}^{(1)} + C_{ub}^{(8)} Q_{ub}^{(8)} + C_{tb}^{(1)} Q_{tb}^{(1)} + C_{tb}^{(8)} Q_{tb}^{(8)} \\ & + [C_{utbd}^{(1)} Q_{utbd}^{(1)} + C_{utbd}^{(8)} Q_{utbd}^{(8)} + \text{h.c.}] \\ & + \begin{cases} \sum_{p=1}^3 (C_{eu})_{pp} Q_{eu,pp} + (C_{ed})_{pp} Q_{ed,pp} + (C_{et})_{pp} Q_{et,pp} + (C_{eb})_{pp} Q_{eb,pp} \\ + \sum_{prst \in P_{ee}} (C_{ee})_{prst} Q_{ee,prst}, & U(1)_{l+e}^3 [\text{top}] \\ \sum_{p,r=1}^3 C_{eu} \delta_{pr} Q_{eu,pr} + C_{ed} \delta_{pr} Q_{ed,pr} + C_{et} \delta_{pr} Q_{et,pr} + C_{eb} \delta_{pr} Q_{eb,pr} \\ + \sum_{p,r,s,t=1}^3 C_{ee} \delta_{pr} \delta_{st} Q_{ee,prst}, & U(3)^2 [\text{topU31}] \end{cases} \end{aligned} \quad (3.92)$$

and it depends on 40 (27) real independent parameters in the **top** (**topU31**) case.

Class 8 c. No basis rotation is required in $\mathcal{L}_6^{(8c)}$, and the quark currents are mapped directly. In the lepton sector, Q_{le} admits 1 independent contraction in the $U(3)^2$ case

(neglecting the subleading contribution $\propto Y_l^2$) and 12 in the $U(1)_{l+e}^3$ case. We choose those with indices $prst$ in the set

$$P_{le} = P_{le}^h \cup P_{le}^{nh}, \quad (3.93)$$

$$P_{le}^h = \{1111, 2222, 3333, 1122, 1133, 2233, 2211, 3311, 3322\}, \quad P_{le}^{nh} = \{1221, 1331, 2332\},$$

where the contractions in P_{le}^h are Hermitian and those in P_{le}^{nh} are not. The operators $Q_{qQtu}^{(1),(8)}$, $Q_{qQbd}^{(1),(8)}$ are not Hermitian and therefore the associated Wilson coefficients are complex.

The Lagrangian reads

$$\Lambda^2 \mathcal{L}_6^{(8c)} = C_{qu}^{(1)} Q_{qu}^{(1)} + C_{qt}^{(1)} Q_{qt}^{(1)} + C_{Qu}^{(1)} Q_{Qu}^{(1)} + C_{Qt}^{(1)} Q_{Qt}^{(1)} \quad (3.94)$$

$$+ C_{qu}^{(8)} Q_{qu}^{(8)} + C_{qt}^{(8)} Q_{qt}^{(8)} + C_{Qu}^{(8)} Q_{Qu}^{(8)} + C_{Qt}^{(8)} Q_{Qt}^{(8)}$$

$$+ C_{qd}^{(1)} Q_{qd}^{(1)} + C_{qb}^{(1)} Q_{qb}^{(1)} + C_{Qd}^{(1)} Q_{Qd}^{(1)} + C_{Qb}^{(1)} Q_{Qb}^{(1)}$$

$$+ C_{qd}^{(8)} Q_{qd}^{(8)} + C_{qb}^{(8)} Q_{qb}^{(8)} + C_{Qd}^{(8)} Q_{Qd}^{(8)} + C_{Qb}^{(8)} Q_{Qb}^{(8)}$$

$$+ \left[C_{qQtu}^{(1)} Q_{qQtu}^{(1)} + C_{qQtu}^{(8)} Q_{qQtu}^{(8)} + C_{qQbd}^{(1)} Q_{qQbd}^{(1)} + C_{qQbd}^{(8)} Q_{qQbd}^{(8)} + \text{h.c.} \right]$$

$$+ \begin{cases} \sum_{p=1}^3 (C_{lu})_{pp} Q_{lu,pp} + (C_{ld})_{pp} Q_{ld,pp} + (C_{lt})_{pp} Q_{lt,pp} + (C_{lb})_{pp} Q_{lb,pp} \\ \quad + (C_{qe})_{pp} Q_{qe,pp} + (C_{Qe})_{pp} Q_{Qe,pp} + \sum_{prst \in P_{le}^h} (C_{le})_{prst} Q_{le,prst} & U(1)_{l+e}^3 [\text{top}] \\ \quad + \left[\sum_{prst \in P_{le}^{nh}} (C_{le})_{prst} Q_{le,prst} + \text{h.c.} \right], \\ \sum_{p,r=1}^3 C_{lu} \delta_{pr} Q_{lu,pr} + C_{ld} \delta_{pr} Q_{ld,pr} + C_{lt} \delta_{pr} Q_{lt,pr} + C_{lb} \delta_{pr} Q_{lb,pr} \\ \quad + C_{qe} \delta_{pr} Q_{qe,pr} + C_{Qe} \delta_{pr} Q_{Qe,pr} + \sum_{p,r,s,t=1}^3 C_{le} \delta_{pr} \delta_{st} Q_{le,prst}, & U(3)^2 [\text{topU31}] \end{cases}$$

and it depends on 54 (31) independent real parameters in the **top** (**topU31**) case.

Class 8 d. Finally, the operators in $\mathcal{L}_6^{(8d)}$ are also mapped directly to the notation with 6 quark fields. $U(2)^3$ invariance requires an insertion of a light Yukawa couplings for each $(\bar{q}u)$ or $(\bar{q}d)$ current and an insertion of Y_l for each $(\bar{l}e)$ current, as indicated in table 2.

This class includes a total of 64 (40) real parameters in the **top** (**topU31**) case:

$$\Lambda^2 \mathcal{L}_6^{(8d)} = C_{quqd}^{(1)} Q_{quqd}^{(1)} + C_{quqd}^{(8)} Q_{quqd}^{(8)} + C_{quqd}^{(1)'} Q_{quqd}^{(1)'} + C_{quqd}^{(8)'} Q_{quqd}^{(8)'} + C_{QtQb}^{(1)} Q_{QtQb}^{(1)} + C_{QtQb}^{(8)} Q_{QtQb}^{(8)} \quad (3.95)$$

$$+ C_{Qtqd}^{(1)} Q_{Qtqd}^{(1)} + C_{quQb}^{(1)} Q_{quQb}^{(1)} + C_{Quqb}^{(1)} Q_{Quqb}^{(1)} + C_{qtQd}^{(1)} Q_{qtQd}^{(1)}$$

$$+ C_{Qtqd}^{(8)} Q_{Qtqd}^{(8)} + C_{quQb}^{(8)} Q_{quQb}^{(8)} + C_{Quqb}^{(8)} Q_{Quqb}^{(8)} + C_{qtQd}^{(8)} Q_{qtQd}^{(8)}$$

$$+ \begin{cases} \sum_{p=1}^3 (C_{lequ}^{(1)})_{pp} Q_{lequ,pp}^{(1)} + (C_{leQt}^{(1)})_{pp} Q_{leQt,pp}^{(1)} + (C_{lequ}^{(3)})_{pp} Q_{lequ,pp}^{(3)} \\ \quad + (C_{leQt}^{(3)})_{pp} Q_{leQt,pp}^{(3)} + (C_{ledq})_{pp} Q_{ledq,pp} + (C_{lebQ})_{pp} Q_{lebQ,pp}, & U(1)_{l+e}^3 [\text{top}] \\ \sum_{p,r=1}^3 C_{lequ}^{(1)} (Y_l^\dagger)_{pr} Q_{lequ,pr}^{(1)} + C_{leQt}^{(1)} (Y_l^\dagger)_{pr} Q_{leQt,pr}^{(1)} + C_{lequ}^{(3)} (Y_l^\dagger)_{pr} Q_{lequ,pr}^{(3)} \\ \quad + C_{leQt}^{(3)} (Y_l^\dagger)_{pr} Q_{leQt,pr}^{(3)} + C_{ledq} (Y_l^\dagger)_{pr} Q_{ledq,pr} + C_{lebQ} (Y_l^\dagger)_{pr} Q_{lebQ,pr}, & U(3)^2 [\text{topU31}] \end{cases}$$

$$+ \text{h.c.}$$

3.5 Comparison with the literature

We conclude this section with a comparison of the parameterizations presented in this section with other recent results in the literature. As a quantitative reference, table 3

	general		U35		MFV		top		topU31	
	all	\mathcal{CP}	all	\mathcal{CP}	all	\mathcal{CP}	all	\mathcal{CP}	all	\mathcal{CP}
$\mathcal{L}_6^{(1)}$	4	2	4	2	2	-	4	2	4	2
$\mathcal{L}_6^{(2,3)}$	3	-	3	-	3	-	3	-	3	-
$\mathcal{L}_6^{(4)}$	8	4	8	4	4	-	8	4	8	4
$\mathcal{L}_6^{(5)}$	54	27	6	3	7	-	14	7	10	5
$\mathcal{L}_6^{(6)}$	144	72	16	8	20	-	36	18	28	14
$\mathcal{L}_6^{(7)}$	81	30	9	1	14	-	21	2	15	2
$\mathcal{L}_6^{(8a)}$	297	126	8	-	10	-	31	-	16	-
$\mathcal{L}_6^{(8b)}$	450	195	9	-	19	-	40	2	27	2
$\mathcal{L}_6^{(8c)}$	648	288	8	-	28	-	54	4	31	4
$\mathcal{L}_6^{(8d)}$	810	405	14	7	13	-	64	32	40	20
tot	2499	1149	85	25	120	-	275	71	182	53

Table 3. Number of independent real parameters in each class of dimension 6 operators, for the 5 flavor structures implemented in SMEFTsim.

summarizes the number of independent real parameters for each class of \mathcal{L}_6 operators and flavor setup.

Compared to previous versions of SMEFTsim [1], the following changes were made:

- the dependence on the CKM matrix in currents involving left-handed down quarks was neglected in the effective operators defined in the **general** and **MFV** versions. It has been restored in version 3.0.
- four parameters corresponding to the real and imaginary parts of $C_{quqd}^{(1)'}$, $C_{quqd}^{(8)'}$ were missing in the **U35** and **MFV** models, and have now been included.
- the **MFV** models have been modified: all Yukawas are now retained in the spurions, instead of only y_t, y_b . Moreover, the Lagrangian is now organized according to a power counting in the quark Yukawas, that led to some flavor-violating terms (eg. $\Delta^u C_{Hud}, \Delta_1^u C_{quqd}^{(1)} \dots$) being dropped, and others (eg. $\Delta^d C_{uH}, C_{qu}^{(1)'}$...) being added.
- versions **top** and **topU31** are new in version 3.0.

The **U35** and **MFV** models can be compared, for instance, to the $U(3)^5$ spurion analyses presented in refs. [66, 75]. Ref. [75] contains an exhaustive classification of all the flavor spurions associated with SM fermion currents in the presence of a $U(3)^5$ symmetry. In their notation, S^u, S^d correspond to Δ_U, Δ_D respectively, while both S^{qu}, S^{qd} are mapped to Δ_Q . The structure $Y_u Y_d^\dagger$ corresponds to Δ_{UD} and, since we only retain linear insertions of $Y_l, \Delta_L = \Delta_E = 0$ in SMEFTsim. Any other spurion leads to baryon and/or lepton number non-conservation, and therefore does not have an equivalent in the Lagrangian considered here.

Ref. [66] presented a detailed classification of all the $U(3)^5$ and $U(2)^5$ invariant structures in the Warsaw basis. In the $U(3)^5$ case, their results can be directly compared with the parameterizations of the **U35** and **MFV** models in **SMEFTsim**, while the $U(2)^3$ case can be compared (in the quark sector) to the **top** and **topU31** models. We find complete agreement in the characterization of the structures, and the operator countings are consistent once a few differences in the organization of the invariants are taken into account:

- the $U(3)^5$ and $U(2)^5$ Lagrangians in ref. [66] are organized according to a power counting in the Yukawas, while for the **U35**, **top** and **topU31** models in **SMEFTsim** we simply choose to retain the leading invariant for each operator in the Warsaw basis.
- in the **MFV** models we retain terms up to order $(Y_l^1, (Y_d + Y_u)^3)$. This choice is different from the power counting in ref. [66], that truncates at $(Y_l^1, Y_d^1 Y_u^2)$.
- the Lagrangian of the **MFV** models includes spurions $\propto Y_d^2$, that were neglected in ref. [66].
- in the $U(2)$ case, different symmetries were chosen for the lepton sector: $U(2)^3$ in ref. [66] vs $U(1)_{l+e}^3$ and $U(3)^2$ in the **top** and **topU31** models.

The structure of the **top** and **topU31** versions builds upon those of refs. [7, 9, 36]. The main difference compared to these works is that in **SMEFTsim** the parameterization has been systematically extended to all operators of the Warsaw basis, including at the same time CP violating terms, interactions that do not involve the top quark, and spurion insertions of the light quark Yukawas.

4 Input parameters

Once the kinetic terms have been canonically normalized and the flavor structure has been fixed, the Lagrangian parameters can be assigned numerical values, with a procedure that is sometimes referred to as “fixing an input parameter scheme” or “finite renormalization”. This section revisits this procedure in the SM and in the SMEFT, using a general formalism that accounts for terms up to arbitrary EFT order. They can be applied to both tree level and loop calculations but, in the latter case, this procedure needs to be combined with the usual renormalization to reabsorb UV divergences. In sections 4.2, 4.3 these formulas are applied to the Warsaw basis case, to recover the known tree-level results, see e.g. [1, 2, 48, 52, 63, 76, 77]. Aspects specific to the NLO case have been discussed in [51, 53, 54, 56, 78, 79].

The Lagrangian parameters are fixed imposing a set of defining conditions that relate them to (pseudo-)observables: for a Lagrangian with N independent parameters $g = \{g_1 \dots g_N\}$, $M \geq N$ independent input observables $\mathcal{O} = \{\mathcal{O}_1 \dots \mathcal{O}_M\}$ need to be selected. Computing each \mathcal{O}_n in the theory at a chosen perturbative order, one obtains relations

$$\mathcal{O}_n = F_n^{(0)}(g), \quad n = 1 \dots M, \tag{4.1}$$

where $F_n^{(0)}$ denotes a function of the parameters g . If $M = N$ and (4.1) is an invertible system of equations, the solution

$$g_i = K_i^{(0)}(\mathcal{O}), \quad i = 1 \dots N, \tag{4.2}$$

fixes g_i as a function $K_i^{(0)}$ of \mathcal{O} . The numerical values of the parameters g are then univocally determined by measurements of \mathcal{O} .

In the SM case, one has 19 independent parameters, that we can classify as

$\alpha_s, \bar{\theta},$	QCD	
$g_1, g_W, v, \lambda,$	EW+Higgs	
$y_e, y_\mu, y_\tau, y_u, y_c, y_t, y_d, y_s, y_b,$	Yukawas	(4.3)
$\theta_{12}, \theta_{13}, \theta_{23}, \delta,$	CKM	

where we have introduced the CP-violating θ angle of QCD ($\bar{\theta}$) and the CKM angles and CP phase ($\theta_{12}, \theta_{13}, \theta_{23}$ and δ respectively). The procedure outlined above is most often employed to determine the value of $\bar{\theta}$ and of the EW+Higgs and Yukawa parameters. On the other hand, the determination of the CKM parameters and of α_s usually relies on a large number of observables: in these cases, the system (4.1) is not invertible and the parameters' values are extracted via a global fit.

When transitioning from the SM to the SMEFT, a large number of additional parameters enters the Lagrangian, namely the cutoff Λ and the Wilson coefficients C_α . Fixing their numerical values in terms of measured observables is obviously still an open challenge (and indeed the ultimate goal of the present work), so these quantities are necessarily left free in the Lagrangian. Nevertheless, they play a role in the finite renormalization procedure, because the observables \mathcal{O} employed to fix the SM quantities generically receive contributions from higher dimensional operators. Working order by order in the EFT expansion, the relations in (4.1) are modified into⁹

$$\mathcal{O}_n = F_n^{(0)}(g) + \frac{1}{\Lambda^2} F_n^{(2)}(g, C) + \frac{1}{\Lambda^4} F_n^{(4)}(g, C) + \dots \tag{4.4}$$

where C here generically represents the set of relevant Wilson coefficients, that can be associated to operators of any dimension. In cases where the system of eq. (4.1) can be inverted, (4.4) can also be solved expanding around the SM solution. The result has the general form:

$$g_i = K_i^{(0)}(\mathcal{O}) + \frac{1}{\Lambda^2} K_i^{(2)}(\mathcal{O}, C) + \frac{1}{\Lambda^4} K_i^{(4)}(\mathcal{O}, C) + \dots \tag{4.5}$$

where $K_i^{(0)}(\mathcal{O})$ is the SM solution and the following K terms are SMEFT corrections that depend on the Wilson coefficients. The leading term in the solution (4.5) is defined imposing that the SM relation holds:

$$\mathcal{O}_n \equiv F_n^{(0)}(K^{(0)}(\mathcal{O})). \tag{4.6}$$

⁹Terms suppressed by odd powers of Λ are omitted here, as they typically contribute to B and/or L violating observables, that are not relevant for the extraction of SM parameters. Nevertheless, the results derived in this section directly generalize to the case where these contributions are retained.

The explicit form of the remaining K terms is found inserting eq. (4.5) into (4.4), expanding in Λ and requiring that SMEFT corrections cancel order by order in the resulting expression. Iteratively, one finds

$$K_i^{(2)} = -(J^{-1})_{in} F_n^{(2)}, \quad (4.7)$$

$$K_i^{(4)} = -(J^{-1})_{in} \left[F_n^{(4)} + \frac{\partial F_n^{(2)}}{\partial g_k} K_k^{(2)} + \frac{1}{2} \frac{\partial^2 F_n^{(0)}}{\partial g_k \partial g_j} K_k^{(2)} K_j^{(2)} \right], \quad (4.8)$$

\vdots

$$K_i^{(d)} = -(J^{-1})_{in} \left[F_n^{(d)} + \sum_{\substack{m < d \\ m+d_1+\dots+d_D=d}} \frac{1}{D!} \frac{\partial^D F_n^{(m)}}{\partial g_{i_1} \dots \partial g_{i_D}} K_{i_1}^{(d_1)} \dots K_{i_D}^{(d_D)} \right], \quad (4.9)$$

where $(J^{-1})_{in} = \partial g_i / \partial F_n^{(0)}$ is the inverse of the Jacobian matrix

$$J_{ni} = \frac{\partial F_n^{(0)}}{\partial g_i}, \quad (4.10)$$

and in eq. (4.9) the sum runs over all possible terms with $m < d$ and such that $m + d_1 + \dots + d_D = d$. All functions and derivatives appearing explicitly in eqs. (4.7)–(4.9) are evaluated at the SM solution for the parameters $g \equiv K^{(0)}(\mathcal{O})$ and the indices $n, k, j, i_1 \dots i_D$ are implicitly contracted internally and summed over.

A generic *predicted* observable \mathcal{P} inherits a dependence on the corrections $F_n^{(d \geq 2)}$ to the input quantities. Analogous to \mathcal{O} , \mathcal{P} will have the generic form

$$\mathcal{P} = P^{(0)}(g) + \frac{1}{\Lambda^2} P^{(2)}(g, C) + \frac{1}{\Lambda^4} P^{(4)}(g, C) + \dots \quad (4.11)$$

where $P^{(0)}$ is the SM expression and $P^{(d \geq 2)}$ encode direct EFT contributions to \mathcal{P} , induced by effective operators entering the relevant Feynman diagrams. Calculating \mathcal{P} in the SMEFT starting from input quantities \mathcal{O} means inserting the expressions of g in eq. (4.5) into eq. (4.11). This operation introduces “indirect” EFT contributions, that are a direct consequence of the $F_n^{(d \geq 2)}$ terms in eq. (4.4). The dependence on the latter quantities can be made explicit:

$$\mathcal{P} = P^{(0)} + \frac{1}{\Lambda^2} \left[P^{(2)} - A_n F_n^{(2)} \right] + \frac{1}{\Lambda^4} \left[P^{(4)} - A_n F_n^{(4)} - A_n^{(2)} F_n^{(2)} + B_{mn} F_m^{(2)} F_n^{(2)} \right] + \dots \quad (4.12)$$

where the m, n indices are summed over and, as above, all functions are implicitly evaluated at $g \equiv K^{(0)}(\mathcal{O})$. The coefficients A, B are found via chain differentiation:

$$A_n = \frac{\partial P^{(0)}}{\partial F_n^{(0)}} = \frac{\partial P^{(0)}}{\partial g_i} (J^{-1})_{in}, \quad (4.13)$$

$$A_n^{(2)} = \frac{\partial P^{(2)}}{\partial F_n^{(0)}} - \frac{\partial P^{(0)}}{\partial F_m^{(0)}} \frac{\partial F_m^{(2)}}{\partial F_n^{(0)}} = \frac{\partial P^{(2)}}{\partial g_i} (J^{-1})_{in} - A_m \frac{\partial F_m^{(2)}}{\partial g_k} (J^{-1})_{kn}, \quad (4.14)$$

$$\begin{aligned} B_{mn} &= \frac{1}{2} \frac{\partial^2 P^{(0)}}{\partial F_m^{(0)} \partial F_n^{(0)}} - \frac{1}{2} \frac{\partial P^{(0)}}{\partial F_p^{(0)}} \frac{\partial^2 F_p^{(0)}}{\partial F_m^{(0)} \partial F_n^{(0)}} = \\ &= \frac{1}{2} \frac{\partial^2 P^{(0)}}{\partial g_i \partial g_j} (J^{-1})_{im} (J^{-1})_{jn} - \frac{1}{2} A_p \frac{\partial^2 F_p^{(0)}}{\partial g_i \partial g_j} (J^{-1})_{im} (J^{-1})_{jn}. \end{aligned} \quad (4.15)$$

Here A_n and the first term in $A_n^{(2)}$ account for linear $K^{(2)}$ corrections to g in the $P^{(0)}$ and $P^{(2)}$ function respectively. The first term in B_{mn} contains double $K^{(2)}$ insertions¹⁰ in $P^{(0)}$, while the second terms of $A_n^{(2)}$ and B_{mn} both stem from $K^{(4)}$ contributions in $P^{(0)}$.

The net effect of the finite renormalization procedure is that all the EFT corrections to input measurements are recast into corrections to predicted quantities: if \mathcal{P} is an input observable $\mathcal{P} \equiv \mathcal{O}_q$, all EFT corrections in eq. (4.12) cancel order by order in the EFT. This happens by construction and follows trivially from the defining conditions imposed. It can be checked explicitly: in this case $P^{(d)} = F_q^{(d)}$ and assuming that \mathcal{O} is a set of independent quantities, also $\partial F_q^{(0)}/\partial F_n^{(0)} = \delta_{qn}$ and $\partial^2 F_q^{(0)}/\partial F_m^{(0)}\partial F_n^{(0)} = 0$. This immediately leads to

$$P^{(2)} - A_n F_n^{(2)} = 0, \quad B_{mn} = 0, \quad (4.16)$$

$$A_n^{(2)} = 0, \quad P^{(4)} - A_n F_n^{(4)} = 0. \quad (4.17)$$

Eq. (4.12) provides a dictionary between different input parameter schemes: comparing sets \mathcal{O} and \mathcal{O}' , the difference in the predicted \mathcal{P} is

$$\begin{aligned} \mathcal{P}(\mathcal{O}) - \mathcal{P}(\mathcal{O}') = & \frac{1}{\Lambda^2} \left[-A_n F_n^{(2)} + A_n' F_n^{(2)'} \right] + \frac{1}{\Lambda^4} \left[-A_n F_n^{(4)} + A_n' F_n^{(4)'} + \right. \\ & \left. - A_n^{(2)} F_n^{(2)} + A_n^{(2)'} F_n^{(2)'} + B_{mn} F_m^{(2)} F_n^{(2)} - B_{mn}' F_m^{(2)'} F_n^{(2)'} \right] + \dots \end{aligned} \quad (4.18)$$

which is easily evaluated via eqs. (4.13)–(4.15). This result is consistent with those in the appendix of ref. [63] and in ref. [80].

4.1 Implementation in SMEFTsim

SMEFTsim implements the finite renormalization procedure via replacements of the form¹¹

$$g_i \rightarrow \hat{g}_i + \delta g_i, \quad (4.19)$$

where \hat{g}_i satisfies the SM relation $\hat{g}_i = K_i^{(0)}(\mathcal{O})$ and δg_i encodes all the dependence on the Wilson coefficients. In the `FeynRules` models, these replacements are operated at the Lagrangian level via the lists `redefConst` (applied simultaneously to the redefinitions in eq. (2.11)) and `redefVev`, and the hats are subsequently dropped in the notation. In this way, all the SM parameters appearing explicitly in the final $\mathcal{L}_{\text{SMEFT}}$ are hatted quantities, i.e. they are conveniently defined in the exact same way as in the SM and their numerical value is directly defined by the input observables chosen.

The shifts δg appear explicitly in the interaction terms, and they are responsible for propagating input shifts corrections to the computed processes. By construction, the de-

¹⁰Here “double insertions” refers to any contribution quadratic in the \mathcal{L}_6 coefficients. This includes contributions from the square of a diagram with one EFT insertion, as well as from the interference between SM and EFT diagrams with two EFT vertices, or EFT diagrams with a single interaction $\propto C^2$. The latter generally stem from field or parameter redefinitions in the Lagrangian.

¹¹Ref. [63] used the notation $\bar{g}_i \rightarrow \hat{g}_i + \delta g_i$ from ref. [52]. This is completely equivalent to the one used here, dropping the bars.

pendence on δg themselves is universal, while their expressions in terms of Wilson coefficients are fixed by the input scheme choice:

$$\delta g_i = \frac{1}{\Lambda^2} K_i^{(2)}(\mathcal{O}) = -\frac{1}{\Lambda^2} (J^{-1})_{in} F_n^{(2)}. \quad (4.20)$$

As noted in section 2, because we work at order Λ^{-2} , the replacements of eq. (4.19) only need to be performed on \mathcal{L}_{SM} and only linear terms in δg need to be retained. Moreover, one can replace $v_T \rightarrow \hat{v}$ in the \bar{C}_α notation, eq. (2.4).

This procedure is implemented for parameters listed in the Higgs, EW and Yukawa sectors in (4.3), as described below. Eq. (4.4) makes manifest that the extraction of SM parameters from global fits can become problematic when generalized to the SMEFT. Whenever this set of equations is not invertible, it is not possible to find a simple form for g_i that expands around the SM solution. A consistent treatment of EFT corrections to such input observables would require to extract simultaneously g_i and C_α , which can be very unpractical or even unfeasible, in the presence of blind directions.

In the case of the CKM parameters, this issue has been overcome in ref. [81], where an optimal set of 4 input measurements was proposed, that allows for a treatment of the CKM angles and phase analogous to that of EW parameters. Its implementation is left for future versions of `SMEFTsim`.

The case of α_s poses a bigger challenge. The strong coupling constant can be determined from a particularly vast range of processes [82], and its extraction is often correlated to that of other physical quantities, such as parton distribution functions (PDFs). A proof-of-concept analysis of SMEFT effects on the PDFs determination was presented in ref. [83], that explored the consequences of including four-fermion operators in a fit to deep-inelastic scattering data. Further studies are needed in order to define an optimal strategy for the treatment of SMEFT contributions in this context. For the time being, input shift corrections associated to the determination of α_s are omitted in `SMEFTsim`.

4.2 Higgs and EW sectors

The electroweak sector of the SM contains 4 independent quantities, that can be chosen as $g = \{g_1, g_W, v, \lambda\}$. The 4 (pseudo-)observables needed to fix their values are usually taken in the set

$$\{\alpha_{\text{em}}, G_F, m_Z, m_W, m_h\}.$$

While m_h always needs to be retained in order to fix λ , the choice of the 3 remaining inputs is free, and several combinations have been adopted in the literature. `SMEFTsim` implements the two alternative schemes $\{\alpha_{\text{em}}, m_Z, G_F\}$ and $\{m_W, m_Z, G_F\}$, providing independent UFO models for both.

The fine structure constant $\alpha_{\text{em}}(0)$ is taken to be measured in Thomson scattering,¹² the Fermi constant G_F measured in muon decays $\mu^- \rightarrow e^- \nu_\mu \bar{\nu}_e$, and m_W, m_Z, m_h are

¹²As we work at tree level, only direct SMEFT corrections to Thomson scattering (i.e. to the determination of $\alpha_{\text{em}}(0)$) are included here. The determination of $\alpha_{\text{em}}(m_Z)$ at one loop in the SMEFT is another major open problem, as potential EFT contributions in the running have not been estimated to date. The main challenge in this task is posed by non-perturbative effects, particularly those arising as α_{em} runs through the hadronic resonances region.

$\Delta\kappa_H = \bar{C}_{H\Box} - \frac{\bar{C}_{HD}}{4}$ $\Delta m_Z^2 = \frac{\bar{C}_{HD}}{2} + \frac{2g_1g_W}{g_1^2 + g_W^2} \bar{C}_{HWB}$ $\Delta\alpha_{\text{em}} = -\frac{2g_1g_W}{g_1^2 + g_W^2} \bar{C}_{HWB}$ $\Delta m_h^2 = 2\Delta\kappa_H - \frac{3}{2\lambda} \bar{C}_H$	<p>general</p> <p>U35, MFV, topU31</p> <p>top</p>	$\Delta G_F = (\bar{C}_{Hl}^{(3)})_{11} + (\bar{C}_{Hl}^{(3)})_{22} - (\bar{C}_{ll})_{1221}$ $\Delta G_F = 2\bar{C}_{Hl}^{(3)} - \bar{C}'_{ll}$ $\Delta G_F = \bar{C}_{Hl,11}^{(3)} + \bar{C}_{Hl,22}^{(3)} - \frac{\bar{C}_{ll,1221}}{2}$
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Table 4. Expressions of input parameter shifts and the kinetic correction $\Delta\kappa_H$ (defined in (2.6)) in terms of Wilson coefficients. The left column is common to all flavor versions, while ΔG_F varies as indicated in the right column. We use the notation $\bar{C}_\alpha = C_\alpha(\hat{v}^2/\Lambda^2)$.

defined as the bosons' pole masses, see ref. [1] and references therein for further details. With these definitions, at tree level:¹³

$$\alpha_{\text{em}} = \frac{1}{4\pi} \frac{g_W^2 g_1^2}{g_W^2 + g_1^2} [1 + \Delta\alpha_{\text{em}}], \quad G_F = \frac{1}{\sqrt{2}v_T^2} [1 + \Delta G_F], \quad (4.21)$$

$$m_W^2 = \frac{v_T^2}{4} g_W^2, \quad m_Z^2 = \frac{v_T^2}{4} (g_W^2 + g_1^2) [1 + \Delta m_Z^2], \quad (4.22)$$

$$m_h^2 = 2v_T^2 \lambda [1 + \Delta m_h^2]. \quad (4.23)$$

The Δ quantities are dimensionless and defined in table 4: ΔG_F is inferred computing the muon decay width at tree level in the SMEFT, while the remaining shifts can be read from the relevant Lagrangian terms. In particular, the contributions in \bar{C}_{HWB} to $\Delta\alpha_{\text{em}}$, Δm_Z^2 follow directly from eq. (2.18) and Δm_h^2 follows from eq. (2.8).

The relations (4.21)–(4.23) can be directly mapped to the notation of eq. (4.4): for instance

$$F_{G_F}^{(0)} = \frac{1}{\sqrt{2}v_T^2}, \quad \frac{1}{\Lambda^2} F_{G_F}^{(2)} = F_{G_F}^{(0)} \Delta G_F, \quad (4.24)$$

and analogously for the other observables.

4.2.1 $\{\alpha_{\text{em}}, m_Z, G_F\}$ scheme

Solving 3 of the 4 eqs. in (4.21), (4.22), plus eq. (4.23), gives expressions for the SM parameters of the form of (4.5). Let us choose the input quantities $\mathcal{O}^{(\alpha)} = \{\alpha_{\text{em}}, m_Z^2, G_F, m_h^2\}$.

It is convenient to define the vector of SM parameters as $g = \{g_1^2, g_W^2, v_T^2, \lambda\}$. The SM solutions $\hat{g}_i \equiv K_i^{(0)}(\mathcal{O})$ are then

$$\hat{g}_1^2 = \frac{4\pi\alpha_{\text{em}}}{c_\theta^2}, \quad \hat{g}_W^2 = \frac{4\pi\alpha_{\text{em}}}{s_\theta^2}, \quad \hat{v}^2 = \frac{1}{\sqrt{2}G_F}, \quad \hat{\lambda} = \frac{m_h^2 G_F}{\sqrt{2}}, \quad (4.25)$$

¹³The normalization of ΔG_F has been modified compared to previous SMEFTsim versions in order to homogenize the notation with the remaining shifts.

The second line was evaluated with generic flavor indices for ΔG_F , and it can be easily mapped to other flavor structures with the dictionary in appendix C. Finally, it is worth noting that electromagnetic interactions do not receive any corrections in this scheme:

$$\frac{\delta e}{\hat{e}} = c_\theta^2 \frac{\delta g_1}{\hat{g}_1} + s_\theta^2 \frac{\delta g_W}{\hat{g}_W} + \frac{1}{2} \Delta \alpha_{\text{em}} = 0, \quad (4.35)$$

consistent with α_{em} being an input quantity.

4.2.2 $\{m_W, m_Z, G_F\}$ scheme

Choosing the input observables $\mathcal{O}^{(m_W)} = \{m_W^2, m_Z^2, G_F, m_h^2\}$, the SM expressions $\hat{g}_i \equiv K_i^{(0)}(\mathcal{O})$ for the relevant parameters are

$$\hat{g}_1^2 = 4\sqrt{2}G_F m_Z^2 s_\theta^2, \quad \hat{g}_W^2 = 4\sqrt{2}G_F m_W^2, \quad \hat{v}^2 = \frac{1}{\sqrt{2}G_F}, \quad \hat{\lambda} = \frac{m_h^2 G_F}{\sqrt{2}}, \quad (4.36)$$

with the weak angle defined by

$$s_\theta^2 \equiv \frac{\hat{g}_1^2}{\hat{g}_1^2 + \hat{g}_W^2} = 1 - \frac{m_W^2}{m_Z^2}. \quad (4.37)$$

The Jacobian $J = \partial \mathcal{O}^{(m_W)} / \partial g$ takes the form

$$J = \frac{1}{4} \begin{pmatrix} \hat{v}^2 & \hat{g}_W^2 & & & \\ \hat{v}^2 & \hat{v}^2 & \hat{g}_W^2/c_\theta^2 & & \\ & & -2\sqrt{2}/\hat{v}^4 & & \\ & & 8\hat{\lambda} & 8\hat{v}^2 & \end{pmatrix}, \quad (4.38)$$

and from eq. (4.7) one has

$$g_1^2 = \hat{g}_1^2 \left[1 + 2 \frac{\delta g_1}{\hat{g}_1} \right], \quad \frac{\delta g_1}{\hat{g}_1} = -\frac{1}{2} \left[\Delta G_F + \frac{\Delta m_Z^2}{s_\theta^2} \right], \quad (4.39)$$

$$g_W^2 = \hat{g}_W^2 \left[1 + 2 \frac{\delta g_W}{\hat{g}_W} \right], \quad \frac{\delta g_W}{\hat{g}_W} = -\frac{\Delta G_F}{2}, \quad (4.40)$$

$$v_T^2 = \hat{v}^2 \left[1 + 2 \frac{\delta v}{\hat{v}} \right], \quad \frac{\delta v}{\hat{v}} = \frac{\Delta G_F}{2}, \quad (4.41)$$

$$\lambda = \hat{\lambda} \left[1 - \frac{\delta \lambda}{\hat{\lambda}} \right], \quad \frac{\delta \lambda}{\hat{\lambda}} = -\Delta G_F - \Delta m_h^2. \quad (4.42)$$

With this input scheme choice, α_{em} is now a predicted quantity. From eq. (4.12):

$$\begin{aligned} \alpha_{\text{em}} &= P_{\alpha_{\text{em}}}^{(0)} + \frac{1}{\Lambda^2} \left[\frac{\partial P_{\alpha_{\text{em}}}^{(0)}}{\partial g_i} (J^{-1})_{in} F_n^{(2)} + P_{\alpha_{\text{em}}}^{(2)} \right] \\ &= \frac{1}{4\pi} \frac{\hat{g}_1^2 \hat{g}_W^2}{\hat{g}_1^2 + \hat{g}_W^2} \left[1 - \Delta G_F - \frac{c_\theta^2}{s_\theta^2} \Delta m_Z^2 + \Delta \alpha_{\text{em}} \right] \\ &= \frac{1}{4\pi} \frac{\hat{g}_1^2 \hat{g}_W^2}{\hat{g}_1^2 + \hat{g}_W^2} \left[1 + 2c_\theta^2 \frac{\delta g_1}{\hat{g}_1} + 2s_\theta^2 \frac{\delta g_W}{\hat{g}_W} + \Delta \alpha_{\text{em}} \right] \\ &= \frac{\hat{e}^2}{4\pi} \left[1 + 2 \frac{\delta e}{\hat{e}} \right]. \end{aligned} \quad (4.43)$$

	ΔM_l	ΔM_u		ΔM_d	
general	$\frac{1}{2}\bar{C}_{eH}^\dagger$	$\frac{1}{2}\bar{C}_{uH}^\dagger$		$\frac{1}{2}\bar{C}_{dH}^\dagger V$	
U35	$\frac{1}{2}\bar{C}_{eH}^* Y_l^{(d)}$	$\frac{1}{2}\bar{C}_{uH}^* Y_u^{(d)}$		$\frac{1}{2}\bar{C}_{dH}^* Y_d^{(d)}$	
MFV	$\frac{1}{2}\bar{C}_{eH} Y_l^{(d)}$	$\frac{1}{2}Y_u^{(d)} [\bar{C}_{uH}^{(0)} + (\Delta^u \bar{C}_{uH}) S^{qu} + (\Delta^d \bar{C}_{uH}) S^{qd}]$		$\frac{1}{2}Y_d^{(d)} V^\dagger [\bar{C}_{dH}^{(0)} + (\Delta^u \bar{C}_{dH}) S^{qu} + (\Delta^d \bar{C}_{dH}) S^{qd}] V$	
top	$\frac{1}{2}(\bar{C}_{eH})_{pp}^*$	$\frac{1}{2}\bar{C}_{uH}^* Y_u^{(d)}$	(u, c)	$\frac{1}{2}\bar{C}_{dH}^* Y_d^{(d)}$	(d, s)
		$\frac{1}{2}\bar{C}_{tH}^*$	(t)	$\frac{1}{2}\bar{C}_{bH}^*$	(b)
topU31	$\frac{1}{2}\bar{C}_{eH}^* Y_l^{(d)}$	$\frac{1}{2}\bar{C}_{uH}^* Y_u^{(d)}$	(u, c)	$\frac{1}{2}\bar{C}_{dH}^* Y_d^{(d)}$	(d, s)
		$\frac{1}{2}\bar{C}_{tH}^*$	(t)	$\frac{1}{2}\bar{C}_{bH}^*$	(b)

Table 5. SMEFT corrections to the fermion mass matrices for each flavor assumption, see section 3 for all definitions. All Wilson coefficients are scalar quantities, except in the **general** case, where they are 3×3 matrices. In the **MFV** case all parameters are real. In the **top** and **topU31** cases $\Delta M_{u,d}, Y_{u,d}$ are 2×2 matrices for the first two generations, and the mass term for the t, b quarks are independent. We use the notation $\bar{C}_\alpha = C_\alpha(\hat{v}^2/\Lambda^2)$ and the results are given in the mass basis of the up-quarks and charged leptons.

It can be instructive to write the final form of the Higgs potential, once the input shifts are applied onto eq. (2.8). For both input schemes considered here, the result is

$$\begin{aligned}
 V(H) + \mathcal{L}_6 = & h^2 \hat{\lambda} \hat{v}^2 + h^3 \hat{\lambda} \hat{v} \left[1 - \frac{\Delta G_F}{2} + \Delta \kappa_H - \frac{1}{\hat{\lambda}} \bar{C}_H \right] \\
 & + h^4 \frac{\hat{\lambda}}{4} \left[1 - \Delta G_F + 2\Delta \kappa_H - \frac{6}{\hat{\lambda}} \bar{C}_H \right] - \frac{3}{4} \frac{h^5}{\hat{v}} \bar{C}_H - \frac{1}{8} \frac{h^6}{\hat{v}^2} \bar{C}_H.
 \end{aligned} \tag{4.44}$$

4.3 Yukawa sector

To fix the SM Yukawa couplings, we take fermion masses as input quantities. From the propagators' poles, at tree level, we have

$$M_\psi = \frac{v_T}{\sqrt{2}} \left[Y_\psi^{(d)} - \Delta M_\psi \right], \tag{4.45}$$

with $Y_\psi^{(d)}$ diagonal. In the **top**, **topU31** cases the index ψ runs over $\psi = \{l, u, d, t, b\}$ so that $M_l, Y_l, \Delta M_l$ are 3×3 tensors, $M_{u,d}, Y_{u,d}, \Delta M_{u,d}$ are 2×2 and $M_{t,b}, Y_{t,b}, \Delta M_{t,b}$ are scalar quantities. In the other flavor setups $\psi = \{l, u, d\}$ and all quantities are 3×3 matrices. The SMEFT corrections ΔM_ψ are given in table 5 for each flavor assumption. The SM solutions are

$$\hat{Y}_\psi^{(d)} = \frac{\sqrt{2}}{\hat{v}} M_\psi, \tag{4.46}$$

and the shifts δY_ψ have the form

$$Y_\psi^{(d)} \rightarrow \hat{Y}_\psi^{(d)} + \delta Y_\psi^{(d)}, \quad \delta Y_\psi^{(d)} = -\frac{\Delta G_F}{2} \hat{Y}_\psi^{(d)} + \Delta M_\psi, \tag{4.47}$$

where ΔG_F enters via eq. (4.41) and ΔM_ψ is non-diagonal and non-Hermitian in general. The expressions (4.46), (4.47) can be easily generalized to setups where M_ψ is not diagonal, by applying the appropriate flavor rotations to both sides of the equations.

The net effect of the finite renormalization procedure is that ΔM_ψ corrections to the fermion mass terms are recast into corrections to the $h\bar{\psi}\psi$ couplings. In unitary gauge, the Lagrangian resulting from the replacements (2.7), (4.47) is

$$\begin{aligned} \mathcal{L}_{\text{Yukawa}} + \mathcal{L}_6^{(5)} = & -\frac{\bar{d}_{R,p} d_{L,r}}{\sqrt{2}} \left[\hat{v} \hat{Y}_d^{(d)} + h \hat{Y}_d^{(d)} \left(1 - \frac{\Delta G_F}{2} + \Delta \kappa_H \right) - \left(2h + \frac{3h^2}{\hat{v}} + \frac{h^3}{\hat{v}^2} \right) \Delta M_d \right]_{pr} \\ & -\frac{\bar{u}_{R,p} u_{L,r}}{\sqrt{2}} \left[\hat{v} \hat{Y}_u^{(d)} + h \hat{Y}_u^{(d)} \left(1 - \frac{\Delta G_F}{2} + \Delta \kappa_H \right) - \left(2h + \frac{3h^2}{\hat{v}} + \frac{h^3}{\hat{v}^2} \right) \Delta M_u \right]_{pr} \\ & -\frac{\bar{e}_{R,p} e_{L,r}}{\sqrt{2}} \left[\hat{v} \hat{Y}_l^{(d)} + h \hat{Y}_l^{(d)} \left(1 - \frac{\Delta G_F}{2} + \Delta \kappa_H \right) - \left(2h + \frac{3h^2}{\hat{v}} + \frac{h^3}{\hat{v}^2} \right) \Delta M_l \right]_{pr} \\ & + \text{h.c.} \end{aligned} \quad (4.48)$$

In the top, topU31 models analogous terms with t, b quarks are also present.

In the `FeynRules` implementation, the common shifts $\delta v, \delta \lambda, \delta Y_\psi$ are automatically replaced with the corresponding expressions in terms of Wilson coefficients. On the other hand, the dependence on the EW shifts $\delta g_1, \delta g_W$ is left explicit in the Lagrangian, as it is identical for all EW input schemes. Once an inputs set is selected, these shifts can be traded for Wilson coefficients expressions: in `Mathematica` this is done via the replacement lists `alphaShifts` or `MwShifts`. In the `UFO` models all shifts are replaced with the Wilson coefficient expressions.

5 SM loop-generated Higgs interactions

Because `SMEFTsim` is designed as a tree-level model, it cannot reproduce processes that only occur at 1-loop. In fact, estimating SMEFT corrections to observables that are genuinely loop-generated both in the SM and at $d = 6$ level is beyond the scope of `SMEFTsim`. Nevertheless, there are cases where a 1-loop SM processes receives *tree* \mathcal{L}_6 corrections. This notably happens in a few relevant Higgs production and decay channels.

In order to enable an estimate of interference terms between \mathcal{L}_6 and SM diagrams for the processes $gg \rightarrow h, h \rightarrow \gamma\gamma, h \rightarrow Z\gamma$, `SMEFTsim` implements effective SM interactions obtained in the large m_t limit. This formally corresponds to matching the SM onto an EFT (we will refer to this as “top-EFT”) where the top quark has been integrated out. The advantage of this approach is that the top loops are effectively reduced to point vertices that can be inserted in tree diagrams. The obvious caveat is that the top-EFT is only valid in a limited kinematic region, as discussed below.

`SMEFTsim 3.0` contains $h\gamma\gamma, hZ\gamma, hgg, hggg$ and $hgggg$ interactions with contributions up to $\mathcal{O}(m_t^{-2})$, i.e. $d = 7$ in the top-EFT. The implemented Lagrangian is:

$$\mathcal{L}_{\text{SMhloop}} = g_{H\gamma\gamma} \mathcal{O}_{\gamma\gamma} + g_{HZ\gamma} \mathcal{O}_{Z\gamma} + g_{Hgg}^{(1)} \mathcal{O}_{gg}^{(1)} + \frac{1}{m_t^2} \sum_{i=2}^5 g_{Hgg}^{(i)} \mathcal{O}_{gg}^{(i)}, \quad (5.1)$$

where

$$\mathcal{O}_{\gamma\gamma} = A_{\mu\nu} A^{\mu\nu} \frac{h}{v}, \quad \mathcal{O}_{z\gamma} = Z_{\mu\nu} A^{\mu\nu} \frac{h}{v}, \quad \mathcal{O}_{gg}^{(1)} = G_{\mu\nu}^a G^{a\mu\nu} \frac{h}{v}, \quad (5.2)$$

$$\mathcal{O}_{gg}^{(2)} = D_\sigma G_{\mu\nu}^a D^\sigma G^{a\mu\nu} \frac{h}{v}, \quad \mathcal{O}_{gg}^{(3)} = f_{abc} G_\mu^{a\nu} G_\nu^{b\sigma} G_\sigma^{c\mu} \frac{h}{v}, \quad (5.3)$$

$$\mathcal{O}_{gg}^{(4)} = D^\mu G_{\mu\nu}^a D_\sigma G^{a\sigma\nu} \frac{h}{v}, \quad \mathcal{O}_{gg}^{(5)} = G^{a\mu\nu} D_\nu D^\sigma G_{\sigma\mu}^a \frac{h}{v}. \quad (5.4)$$

The corresponding g coefficients are fixed via a 1-loop matching procedure of the SM onto the top-EFT. For the $h\gamma\gamma$ and $hZ\gamma$ interactions we use the results from refs. [84–87], that include loops of both top quarks and W bosons:

$$g_{H\gamma\gamma} = \frac{e^2}{8\pi^2} \left[I_w \left(\frac{m_h^2}{4m_W^2} \right) + \frac{4}{3} I_f \left(\frac{m_h^2}{4m_t^2}, 0 \right) \right], \quad (5.5)$$

$$g_{HZ\gamma} = \frac{e^2}{4\pi^2} \left[t_\theta I_w^Z \left(\frac{m_h^2}{4m_W^2}, \frac{m_Z^2}{4m_W^2} \right) + \left(\frac{1}{2} - \frac{4}{3} s_\theta^2 \right) I_f \left(\frac{m_h^2}{4m_t^2}, \frac{m_Z^2}{4m_t^2} \right) \right]. \quad (5.6)$$

The loop functions I_f, I_w, I_w^Z are evaluated in the limit where the Higgs boson is on-shell and higher order corrections are simply obtained via Taylor-expansion, retaining terms up to $\mathcal{O}(m_t^{-2}m_W^{-6})$:

$$I_f(a, b) = \int_0^1 \int_0^{1-x} \frac{1 - 4xy}{1 - 4(a-b)xy - 4by(1-y)} dy dx \quad (5.7)$$

$$= \frac{1}{3} + \frac{7a}{90} + \frac{11b}{90} + \mathcal{O}(a^2, b^2), \quad (5.8)$$

$$I_w(a) = \int_0^1 \int_0^{1-x} \frac{-4 + 6xy + 4axy}{1 - 4axy} dy dx \quad (5.9)$$

$$= -\frac{7}{4} - \frac{11a}{30} - \frac{19a^2}{105} - \frac{58a^3}{525} + \mathcal{O}(a^4), \quad (5.10)$$

$$I_w^Z(a, b) = \frac{1}{t_\theta^2} \int_0^1 \int_0^{1-x} \frac{(5 - t_\theta^2 + 2a(1 - t_\theta^2))xy + t_\theta^2 - 3}{1 - 4(a-b)xy - 4by(1-y)} dy dx \quad (5.11)$$

$$= 1 + \frac{5 - t_\theta^2}{12} a - b + \frac{t_\theta^2 - 9}{9} ab + \frac{2(5 - t_\theta^2)}{45} a^2 + \frac{56b^2}{45} + \frac{5 - t_\theta^2}{35} a^3 - \frac{12b^3}{7} +$$

$$- \frac{2(7 - t_\theta^2)}{15} a^2 b + \frac{17(13 - t_\theta^2)}{105} ab^2 + \mathcal{O}(a^3, b^3). \quad (5.12)$$

For the Higgs-gluon interactions, the matching has been performed in refs. [88–90] up to dimension 7 in the top expansion:¹⁴

$$g_{Hgg}^{(1)} = \frac{g_s^2}{48\pi^2} + \mathcal{O}(g_s^4), \quad (5.13)$$

$$g_{Hgg}^{(2)} = -\frac{7g_s^2}{2880\pi^2} + \mathcal{O}(g_s^4), \quad (5.14)$$

$$g_{Hgg}^{(3)} = \frac{g_s^3}{240\pi^2} + \mathcal{O}(g_s^5), \quad (5.15)$$

$$g_{Hgg}^{(4)} = \frac{g_s^2}{1440\pi^2} + \mathcal{O}(g_s^4), \quad (5.16)$$

$$g_{Hgg}^{(5)} = \frac{g_s^2}{80\pi^2} + \mathcal{O}(g_s^4). \quad (5.17)$$

¹⁴There is a sign difference in the definition of \mathcal{O}_3 compared to refs. [88, 89]. The sign of C_3 is also affected by the sign in the covariant derivative definition, that was taken with the opposite convention in ref. [90].

Note that the $d = 7$ operators produce interactions with one Higgs and up to 6 gluon legs. While the full gauge-invariant Lagrangian is implemented in the `FeynRules` models, only vertices with up to 4 gluon legs ($hgg, hggg, hgggg$) were exported to the UFOs. The Feynman rules of the $hggggg$ and $hgggggg$ vertices are extremely complex both in the color and Lorentz structures, to the point that their inclusion makes the Monte Carlo event generation computationally challenging. They are available upon request.

5.1 Validity of the approximations used

The Higgs interactions described in this section are implemented to the specific purpose of enabling the simulation of Higgs production and decay processes. In general, these vertices *should not* be inserted into other arbitrary processes. In `MadGraph5_aMC@NLO`, the insertions can be controlled at the diagram generation level via the interaction order `SMHLOOP = 1` that is assigned to all the g couplings in the Lagrangian (5.1), see also section 8.2.

The following limitations should also be kept in mind:

- The implementation relies on the top-EFT formalism, that is only valid when the momentum q flowing through the effective vertex is $q < m_t$. This condition is always fulfilled for $gg \rightarrow h$ with no extra jets, for which the top-EFT reproduces the 1-loop SM cross-section within an accuracy of a few permille. With more complex final states, a validity threshold is present and it can translate differently in terms of measured observables, depending on the process.

Figure 1 shows the relative deviation of the top-EFT predictions from the 1-loop SM results for $d\sigma/dp_T(h)$ in $pp \rightarrow hj$ and $pp \rightarrow hjj$, as obtained at parton level with `MadGraph5_aMC@NLO`, in a 4-flavor scheme and neglecting all electroweak contributions. The SM prediction was obtained generating 100000 events for each $pp \rightarrow hj$ channel and for $qq \rightarrow hqq$, and 50000 events for the remaining $pp \rightarrow hjj$ channels, with the `loop_sm` UFO. The associated PDF and scale uncertainties were estimated with the `MadGraph5_aMC@NLO` functionalities [91], and their combination in quadrature is shown for reference as a grey band. The top-EFT predictions were obtained reweighting the events with `SMEFTsim`. The lines in color compare three different implementations of the top-EFT: including all operators up to $d = 7$ (red), including only the $d = 5$ operator $\mathcal{O}_{gg}^{(1)}$ (blue) and including only the ggh vertex as in the previous `SMEFTsim` versions (orange), see the next subsection. The statistical uncertainty associated to each line is shown as a colored band surrounding the solid curves. Because the statistical errors associated to the reweighted histogram and to the original one are fully correlated, in most cases, the uncertainty on their ratio cancels and it is not visible on the plot. Uncertainties due to the reweighting procedure itself have been neglected, in the absence of a prescription for their estimation.

For both $pp \rightarrow hj$ and $pp \rightarrow hjj$, the total cross section is dominated by gg - and qq -initiated channel, for which the $m_t \rightarrow \infty$ approximation breaks down roughly at $p_T(h) \simeq 250$ GeV [89, 90, 92–95]. Within the top-EFT validity regime, the $d = 7$ implementation reproduces the SM 1-loop result within an accuracy of few %, see also ref. [90]. The large m_t approximation fails most significantly in qq -initiated

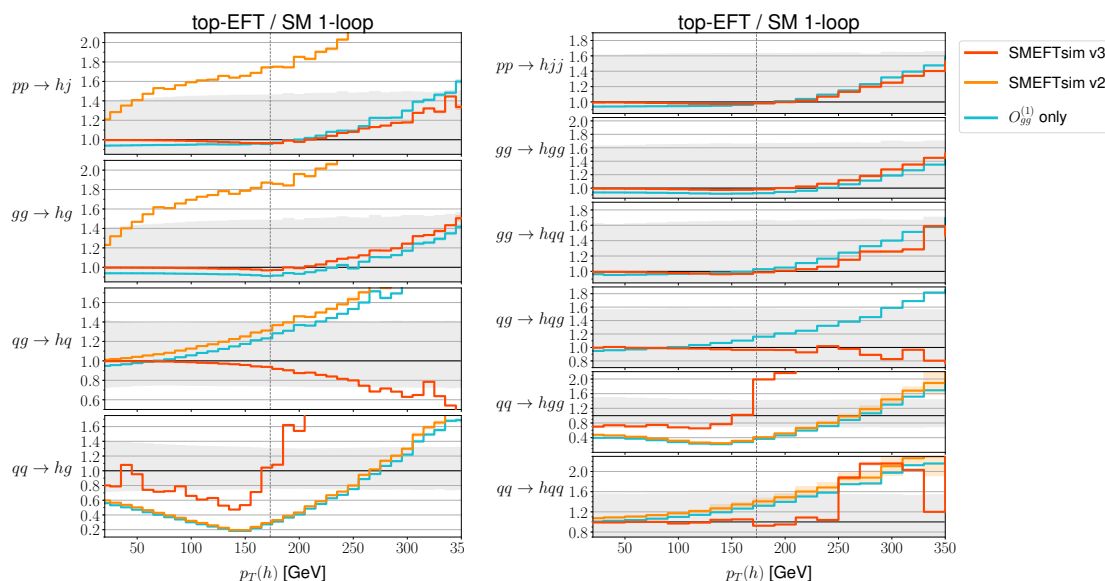


Figure 1. Top-EFT predictions for $d\sigma/dp_T(h)$ in $pp \rightarrow hj$ (left) and $pp \rightarrow hjj$ (right), normalized to the SM 1-loop results, see text for the calculation details. In both figures, the first panel shows the combined result, while the lower ones give the breakdown into the contributing channels. Note that here q generically denotes a light quark or antiquark. The statistical uncertainties are plotted in color, and they are not visible in most cases. For reference, the gray bands show the systematic uncertainty on the SM event generation and the vertical dashed line marks $p_T = m_t$, where the top-EFT is expected to break down. SMEFTsim v2 curves for $pp \rightarrow hjj$, $gg \rightarrow hgg$, $gg \rightarrow hqq$, $qq \rightarrow hqq$ lie above the plotted range.

processes that, nevertheless, give a negligible contribution to the total cross section. This behavior is due to the quarks' PDFs preferring significantly larger x compared to the gluon one, which leads to large \hat{s} contributions being suppressed for gg and qq initial states, but not for qq [94].

- The operators $O_{Hgg}^{(i)}$ form a complete basis up to $\mathcal{O}(m_t^{-2})$ for Higgs interactions with up to 4 gluons [88, 96]. This means that, within the regime of validity of the top-EFT, SMEFTsim can reproduce 1-loop SM Higgs production in gluon fusion with up to 2 jets.

Processes $gg \rightarrow h + nj$ with $n \geq 3$ cannot be fully reproduced with SMEFTsim, even with the inclusion of $hggggg$, $hgggggg$ vertices, because a complete matching to $\mathcal{O}(m_t^{-2})$ onto these vertices would require $d = 9$ top-EFT operators.

- In addition to the validity of the large m_t approximation, the implementation of $h\gamma\gamma$, $hZ\gamma$ assumes an on-shell Higgs in the parameterization of the loop function.

5.2 Comparison to previous versions of SMEFTsim

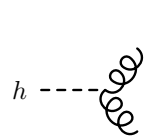
Previous SMEFTsim versions only implemented the hgg , $h\gamma\gamma$ and $hZ\gamma$ vertices, while interactions with higher numbers of gluons were omitted. In version 3.0, all the vertices induced

by the operator $\mathcal{O}_{Hgg}^{(1)}$ and vertices with up to 5 legs (4 gluons) from $\mathcal{O}_{Hgg}^{(2,3,4,5)}$ are included.

Moreover, the hgg interaction was previously parameterized in the on-shell Higgs limit, analogously to $h\gamma\gamma$, $hZ\gamma$, via a coupling [86, 87, 97]

$$g_{Hgg} = \frac{g_s^2}{16\pi^2} I_f \left(\frac{m_h^2}{4m_t^2}, 0 \right),$$

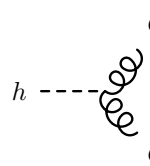
and the loop functions I_f , I_w , I_w^Z were expanded up to $\mathcal{O}(m_t^{-6}m_W^{-6})$, which is formally equivalent to a matching up to $d = 11$ in the top-EFT for an on-shell Higgs boson. In version 3.0 this is replaced by a consistent matching up to $\mathcal{O}(m_t^{-2}m_W^{-6})$ that does not rely on the on-shell assumption. The two parameterizations are completely equivalent up to $\mathcal{O}(m_t^{-4})$ corrections for on-shell Higgs production with no extra jets: the Feynman rule of the ggh interaction is



$$= \frac{4i}{v} [p_1^\nu p_2^\mu - \eta^{\mu\nu} p_1 \cdot p_2] \left[g_{Hgg}^{(1)} - \frac{p_1 \cdot p_2}{m_t^2} g_{Hgg}^{(2)} + \frac{p_1^2 + p_2^2}{4m_t^2} g_{Hgg}^{(5)} \right] +$$

$$- \frac{2i}{v} \frac{p_1^\mu p_1^\nu p_2^2 + p_2^\mu p_2^\nu p_1^2 - \eta^{\mu\nu} p_1^2 p_2^2 - p_1^\mu p_2^\nu p_1 \cdot p_2}{m_t^2} g_{Hgg}^{(4)}, \quad (5.18)$$

with the momenta p_1 , p_2 taken to be incoming. In the limit $p_{1,2}^2 = 0$, $p_h^2 = (p_1 + p_2)^2 = m_h^2$



$$\xrightarrow{\text{on-shell}} \frac{4i}{v} \left[p_1^\nu p_2^\mu - \eta^{\mu\nu} \frac{m_h^2}{2} \right] \left[g_{Hgg}^{(1)} - \frac{m_h^2}{2m_t^2} g_{Hgg}^{(2)} \right] + \frac{i}{v} \frac{m_h^2}{m_t^2} p_1^\mu p_2^\nu g_{Hgg}^{(4)}.$$

(5.19)

The last term vanishes for external gluons and, using eqs. (5.13), (5.14),

$$g_{Hgg}^{(1)} - \frac{m_h^2}{2m_t^2} g_{Hgg}^{(2)} = \frac{g_s^2}{16\pi^2} \left[\frac{1}{3} + \frac{7}{90} \frac{m_h^2}{4m_t^2} \right]. \quad (5.20)$$

The terms in brackets reproduce the expansion of the top loop I_f (eq. (5.8)) up to $\mathcal{O}(m_t^{-2})$.

6 Propagator corrections

Mass terms and decay widths of the SM particles generally receive corrections from \mathcal{L}_6 operators. In order to compute amplitudes consistently at $\mathcal{O}(\Lambda^{-2})$, these corrections need to be included in the propagators.

In unitary gauge the propagator of a generic unstable vector V , scalar S or fermion ψ has the form

$$P_V^{\mu\nu} = \frac{i}{q^2 - m_V^2 + im_V\Gamma_V} \left(-\eta^{\mu\nu} + \frac{q^\mu q^\nu}{m_V^2} \right), \quad (6.1)$$

$$P_S = \frac{i}{q^2 - m_S^2 + im_S\Gamma_S}, \quad (6.2)$$

$$P_\psi = \frac{i(\not{q} + m_\psi)}{q^2 - m_\psi^2 + im_\psi\Gamma_\psi}. \quad (6.3)$$

In the SMEFT we can write, for each particle,

$$m = m^{\text{SM}} + \delta m, \quad \Gamma = \Gamma^{\text{SM}} + \delta \Gamma, \quad (6.4)$$

where the shifts δm , $\delta \Gamma$ collect all the contributions from $d \geq 6$ operators. The corresponding propagator expressions be expanded to linear order in the shifts [98]

$$P_V^{\mu\nu} = P_V^{\mu\nu, \text{SM}} + \Delta P_V^{\mu\nu}, \quad P_S = P_S^{\text{SM}} + \Delta P_S, \quad P_\psi = P_\psi^{\text{SM}} + \Delta P_\psi, \quad (6.5)$$

where the expressions for P^{SM} are given by eqs. (6.1)–(6.3) replacing $m \rightarrow m^{\text{SM}}$, $\Gamma \rightarrow \Gamma^{\text{SM}}$. The corrections read

$$\Delta P_V^{\mu\nu} = -P_V^{\mu\nu, \text{SM}} \frac{im_V^{\text{SM}}}{D_V(q^2)} \left[\delta \Gamma_V + 2i \left(1 - \frac{i\Gamma_V^{\text{SM}}}{2m_V^{\text{SM}}} \right) \delta m_V \right] - 2i \frac{q^\mu q^\nu}{D_V(q^2)} \frac{\delta m_V}{(m_V^{\text{SM}})^3}, \quad (6.6)$$

$$\Delta P_S = -P_S^{\text{SM}} \frac{im_S^{\text{SM}}}{D_S(q^2)} \left[\delta \Gamma_S + 2i \left(1 - \frac{i\Gamma_S^{\text{SM}}}{2m_S^{\text{SM}}} \right) \delta m_S \right], \quad (6.7)$$

$$\Delta P_\psi = -P_\psi^{\text{SM}} \frac{im_\psi^{\text{SM}}}{D_\psi(q^2)} \left[\delta \Gamma_\psi + 2i \left(1 - \frac{i\Gamma_\psi^{\text{SM}}}{2m_\psi^{\text{SM}}} \right) \delta m_\psi \right] + \frac{i\delta m_\psi}{D_\psi(q^2)}, \quad (6.8)$$

with the shorthand notation

$$D(q^2) = q^2 - (m^{\text{SM}})^2 + im^{\text{SM}} \Gamma^{\text{SM}}. \quad (6.9)$$

Note that, since $\Delta P \propto D(q^2)^{-1}$, propagator corrections are expected to be relevant in the on-shell kinematic region and suppressed when the particle is largely off-shell. The dominant contributions are therefore approximated by the on-shell expressions:¹⁵

$$\Delta P_V^{\mu\nu}|_{q^2=m^2} = P_V^{\mu\nu}|_{q^2=m^2} \left[-\frac{\delta \Gamma_V}{\Gamma_V^{\text{SM}}} - \left(1 + \frac{2im_V^{\text{SM}}}{\Gamma_V^{\text{SM}}} \right) \frac{\delta m_V}{m_V^{\text{SM}}} \right], \quad (6.10)$$

$$\Delta P_S|_{q^2=m^2} = P_S|_{q^2=m^2} \left[-\frac{\delta \Gamma_S}{\Gamma_S^{\text{SM}}} - \left(1 + \frac{2im_S^{\text{SM}}}{\Gamma_S^{\text{SM}}} \right) \frac{\delta m_S}{m_S^{\text{SM}}} \right], \quad (6.11)$$

$$\Delta P_\psi|_{q^2=m^2} = P_\psi|_{q^2=m^2} \left[-\frac{\delta \Gamma_\psi}{\Gamma_\psi^{\text{SM}}} - \left(1 + \frac{2im_\psi^{\text{SM}}}{\Gamma_\psi^{\text{SM}}} \right) \frac{\delta m_\psi}{m_\psi^{\text{SM}}} \right] + \frac{\delta m_\psi}{m_\psi^{\text{SM}} \Gamma_\psi^{\text{SM}}}, \quad (6.12)$$

with

$$P_V^{\mu\nu}|_{q^2=m^2} = \frac{-\eta^{\mu\nu}}{m_V \Gamma_V}, \quad P_S|_{q^2=m^2} = \frac{1}{m_S \Gamma_S}, \quad P_\psi|_{q^2=m^2} = \frac{\not{q} + m_\psi}{m_\psi \Gamma_\psi}. \quad (6.13)$$

6.1 Implementation in SMEFTsim

SMEFTsim 3.0 implements propagator corrections for the Z, W, h bosons and for the top quark. The user has two alternative options for including them in SMEFT predictions:

- (a) using the linearized propagator expressions of eqs. (6.5)–(6.8). In this case the pole of the propagator remains located at m^{SM} , and the dependence on the Wilson coefficients, stemming both from δm and $\delta \Gamma$, is linear at the amplitude level. This option is selected fixing `linearPropCorrections = 1` (or any value $\neq 0$) in the `param_card`.

¹⁵Longitudinal contributions for vector bosons were neglected here.

- (b) using the propagator expressions in eqs. (6.1)–(6.3), with shifted masses. In this case the pole of the propagator is located at $m = m^{\text{SM}} + \delta m$ while width corrections are entirely dropped.¹⁶ The dependence on the Wilson coefficients is generally non-linear, as contributions $\propto 1/C_\alpha$ are induced in the amplitude. This is the default option and it's selected with `linearPropCorrections = 0`.

While option (a) is recommended for consistency of the EFT expansion, we caution the user that the linearization can be problematic, particularly in the presence of *mass* corrections. Formally, expanding around the complex pole of the propagator is not a gauge-invariant operation [99–101]. Numerically, significantly large discrepancies between methods (a) and (b) can occur, as illustrated in figure 2 (left) for the case of the W boson. Using linearized propagators leads to sizeable numerical distortions already for $\delta m_W/m_W^{\text{SM}} \gtrsim \text{few } \%$, as the dashed curves show. For comparison, the linear approximation works very well for width corrections up to $\mathcal{O}(10\%)$, see figure 2 (right). This is partially due to the W boson being narrow in the SM. Adopting a $\{m_W, m_Z, G_F\}$ input scheme is convenient in this respect, because mass correction effects are entirely avoided. For a more general discussion of the theoretical advantages of this scheme choice, see e.g. refs. [1, 2, 51, 53, 63, 78, 98, 102].

Option (a) is implemented in `SMEFTsim` following the method outlined in ref. [103], contribution 12:¹⁷ four dummy fields Z', W', h', t' (`Z1, W1, H1, t1`) are introduced, with propagators $\Delta P_Z^{\mu\nu}, \Delta P_W^{\mu\nu}, \Delta P_h, \Delta P_t$. The numerical values assigned to $m^{\text{SM}}, \Gamma^{\text{SM}}$ appearing in these expressions are those of the nominal mass and widths of the dummy fields. The latter, in turn, are internal parameters and defined as equal to the SM masses and widths of the corresponding dynamical particles. The $\delta m, \delta \Gamma$ parameters are also defined as internal parameters, function of the relevant SM couplings and Wilson coefficients.

The dummy states have the same SM interactions as Z, W, h, t and do not enter $d = 6$ operators: `SMEFTsim` contains copies of all SM vertices, with one or more of the standard fields replaced by its dummy counterpart. Vertices with n dummy legs are proportional to n powers of a flag parameter `propCorr` and have interaction order `NPprop = n`. The numerical value of `propCorr` is set to 0(1) if `linearPropCorrections = 0` (a non-zero value).

In this way, for instance, linearized Z -propagator corrections to $pp \rightarrow \mu^+ \mu^-$ can be estimated computing the $pp \rightarrow Z \rightarrow \mu^+ \mu^-$ and $pp \rightarrow Z' \rightarrow \mu^+ \mu^-$ amplitudes, and using the interaction order `NPprop` to isolate the pure SM/interference/quadratic contributions as detailed in section 8.3.

Note that linearized propagator corrections are available only in the UFO models, as the propagators are modified directly in the `propagators.py` file and not in `FeynRules`.¹⁸

Mass and width corrections implemented. All the mass and width shifts implemented in `SMEFTsim` are computed to $\mathcal{O}(\Lambda^{-2})$, i.e. linearly in the Wilson coefficients. Be-

¹⁶The implementation of EFT-corrected decay widths has been avoided to prevent potential conflicts with the treatment of widths in Monte Carlo generators.

¹⁷I thank O. Mattelaer for pointing me to this reference.

¹⁸The expressions in `propagators.py` differ by an overall i factor from those in (6.6), (6.7), because an i is conventionally added by `ALOHA` upon parsing the UFO model [104].

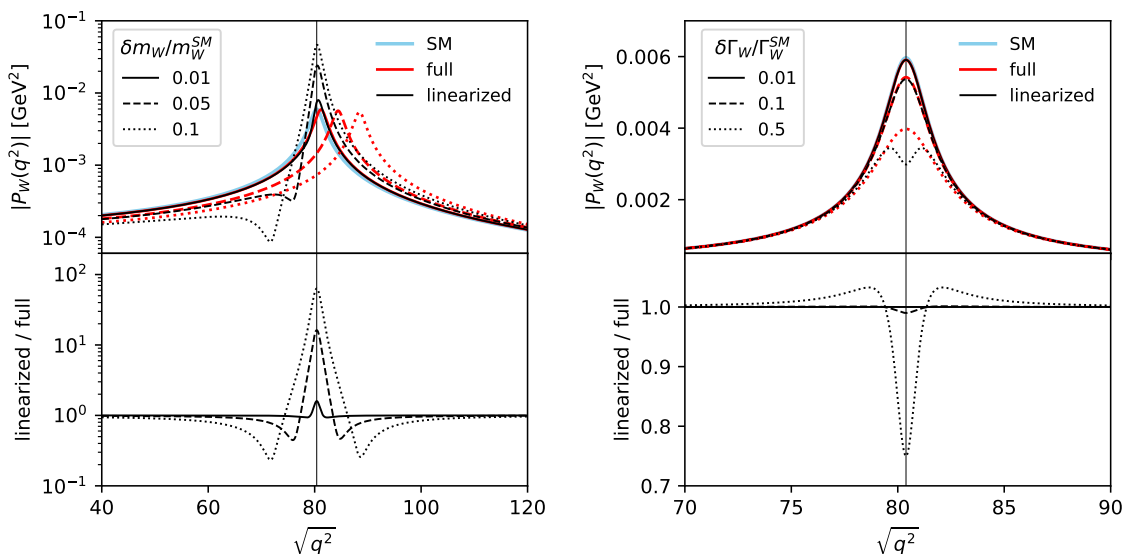


Figure 2. Illustrative comparison of full and linearized W propagators varying δm_W (left) and $\delta \Gamma_W$ (right). The upper panels show the absolute value of the form factor of the transverse propagator $|P_W(q^2)|$ as a function of $\sqrt{q^2}$. Red lines are obtained modifying m_W, Γ_W in the full propagator, eq. (6.1), while black lines are obtained linearizing out $\delta m_W, \delta \Gamma_W$ corrections as in eqs. (6.5), (6.6). The lower panels show the ratio between the two curves. The thin vertical line marks $\sqrt{q^2} = m_W^{SM} = 80.387$ GeV. The light blue curve in the upper plots marks for reference the SM behavior ($\delta \Gamma_W = 0 = \delta m_W$). Solid, dashed and dotted lines correspond to three different sizes of δm_W and $\delta \Gamma_W$, with the same coding in upper and lower panels. In the right panel, the black, red and light blue solid lines are indistinguishable.

cause m_Z, m_h and m_t are taken as input parameters,

$$\delta m_Z \equiv 0, \quad \delta m_h \equiv 0, \quad \delta m_t \equiv 0. \quad (6.14)$$

The δm_W correction is non-vanishing only in the $\{\alpha_{em}, m_Z, G_F\}$ scheme, and the expression was given in eq. (4.34). Decay width corrections for the Z, W bosons and for the top quark are defined as

$$\delta \Gamma_Z = \Gamma_Z^{SM, \text{best}} \left[\frac{\delta \Gamma_Z}{\Gamma_Z^{SM}} \right]_{\text{tree}}, \quad \delta \Gamma_W = \Gamma_W^{SM, \text{best}} \left[\frac{\delta \Gamma_W}{\Gamma_W^{SM}} \right]_{\text{tree}}, \quad \delta \Gamma_t = \Gamma_t^{SM, \text{best}} \left[\frac{\delta \Gamma_t}{\Gamma_t^{SM}} \right]_{\text{tree}}, \quad (6.15)$$

with

$$\Gamma_Z^{SM, \text{best}} = 2.4952 \text{ GeV}, \quad \Gamma_W^{SM, \text{best}} = 2.085 \text{ GeV}, \quad \Gamma_t^{SM, \text{best}} = 1.33 \text{ GeV}, \quad (6.16)$$

the loop-improved SM predictions [82, 105]. These are free parameters in the models, that can be modified by the user. The quantities $\left[\frac{\delta \Gamma}{\Gamma^{SM}} \right]_{\text{tree}}$ are calculated at tree level (both numerator and denominator) using the width computation tools in `FeynRules` [106]. They include all 2-body decays and are extracted in the limit $V_{CKM} = \mathbb{1}$, with all fermion masses set to zero, except those of the b and t quarks. Analytic expressions are given in appendix A.

f	$\text{Br}_{h \rightarrow f}^{SM,\text{best}}$	f	$\text{Br}_{h \rightarrow f}^{SM,\text{best}}$	f	$\text{Br}_{h \rightarrow f}^{SM,\text{best}}$
$\gamma\gamma$	2.27×10^{-3}	$Z\gamma$	1.541×10^{-3}	gg	0.0818
$\bar{b}b$	0.5809	$\bar{c}c$	0.02884	$\tau^+\tau^-$	0.06256
$ll\bar{\nu}\nu$	0.0256	$\bar{u}u\bar{d}d$	0.1097	$l^-\bar{\nu}d\bar{u} + \text{h.c.}$	0.1062

Table 6. Numerical values of the Higgs boson branching ratios employed in the definition of $\delta\Gamma_h$, eq. (6.17). The values for 2-body decays are taken from ref. [107], with $m_h = 125.09$ GeV. The values in the last line are estimated computing the partial widths with **Prophecy4f** 2.0 [108] and normalizing their sum to $\text{Br}_{h \rightarrow 4f}^{SM,\text{best}} = 0.24161$ [107]. They include only charged current contributions and are summed over all allowed flavor combinations.

The correction to the total Higgs width is computed using individual K -factors for each decay channel, as in ref. [57].¹⁹

$$\delta\Gamma_h = \Gamma_h^{SM,\text{best}} \sum_f \text{Br}_{h \rightarrow f}^{SM,\text{best}} \left[\frac{\delta\Gamma_{h \rightarrow f}}{\Gamma_{h \rightarrow f}^{\text{SM}}} \right]_{\text{tree}}, \quad \Gamma_h^{SM,\text{best}} = 4.07 \text{ MeV}, \quad (6.17)$$

with f running over the set $\{\gamma\gamma, Z\gamma, gg, \bar{b}b, \bar{c}c, \tau^+\tau^-\}$ plus the allowed 4-fermion channels. In the **SMEFTsim** implementation, only 4-fermion decays proceeding via charged currents ($h \rightarrow WW^* \rightarrow 4f$) are retained, in order to simplify the analytic expressions. Channels mediated by neutral bosons ($h \rightarrow ZZ^*, Z\gamma^*, \gamma^*\gamma^*, g^*g^* \rightarrow 4f$) give subdominant corrections, that are estimated in a 3 – 5% change to the dependence on $\bar{C}_{HW}, \bar{C}_{HB}, \bar{C}_{HD}$ and a change $\lesssim 1\%$ for the other Wilson coefficients [57].

$\Gamma_h^{SM,\text{best}}$ is a free parameter in the models and can be modified by the user. The best-fit branching ratios, instead, are embedded numerically in the $\delta\Gamma_h$ expressions and cannot be changed. The values employed are reported in table 6. The relative deviations $\left[\frac{\delta\Gamma_{h \rightarrow f}}{\Gamma_{h \rightarrow f}^{\text{SM}}} \right]_{\text{tree}}$ for 2-body decays are computed with the **FeynRules** tools, retaining the full dependence on all the relevant fermion masses and Yukawa couplings. If a given Yukawa coupling y_f is set to zero in the **param_card**, all contributions to $\delta\Gamma_h$ originating from the $h \rightarrow f\bar{f}$ decay channel are dropped. For the $h \rightarrow 4f$ channels we take the analytic results of ref. [57], that neglect *all* fermion masses and quark mixings. Note that the results in ref. [57] were given for the $U(3)^5$ flavor symmetric case, and they have been generalized to the other flavor assumptions in **SMEFTsim**. Full analytic results for SMEFT corrections are reported in appendix A.

7 Usage in Mathematica

The **FeynRules** files in **SMEFTsim** can be imported in **Mathematica** [109] and used to print out analytic expressions for the Feynman rules and Lagrangian terms. A template **Mathematica** notebook is available at the GitHub repository, with examples of usage of the

¹⁹This normalization choice is due to radiative corrections affecting the various channels in significantly different ways. For comparison, in the case of Z, W decays, using individual K -factors leads to variations $\lesssim 2\%$ in the Wilson coefficient dependence compared to an overall rescaling. The top case is trivial, as there is only one relevant decay channel $t \rightarrow bW^+$. In the Higgs case, due to the heterogeneity of the relevant decay processes, the discrepancy between the two normalizations is of order 20–50%.

code in different setups. The functionalities are the standard `FeynRules` ones. However, since `SMEFTsim` is an unusually complex model, some recommendations are in order.

Before importing the model, the user must specify a flavor setup and EW input scheme choice. For instance, after loading `FeynRules`:

```
SetDirectory ["PATH/TO/SMEFTSIM/DIR/"];

Flavor = U35;
Scheme = MwScheme;
LoadModel ["SMEFTsim_main.fr"];
```

Allowed options for `Flavor` are `general`, `U35`, `MFV`, `top`, `topU31`. Allowed options for `Scheme` are `alphaScheme`, `MwScheme`. The loading time varies between flavor assumptions and can take up to a few seconds.²⁰

The `FeynRules` code is split over different files that contain the required operators, parameters and Lagrangian definitions. The implementation is such that only the objects matching the selected flavor structure and EW input scheme are defined upon loading. In all models, the following Lagrangians are defined (definitions were given in section 1):

- `LGauge` = $\mathcal{L}_{\text{gauge}}$. Contains the SM terms plus the linearized SMEFT corrections due to field redefinitions and input parameter shifts.
- `LGaugeP`. Same as `LGauge`, but with at least one W or Z boson replaced with the corresponding dummy field `W1`, `Z1`.
- `LHiggs` = $\mathcal{L}_{\text{Higgs}}$. Contains the SM terms plus the linearized SMEFT corrections due to field redefinitions and input parameter shifts.
- `LHiggsP`. Same as `LHiggs`, but with at least one W , Z or Higgs boson replaced with the corresponding dummy field `W1`, `Z1`, `H1`.
- `LFermions` = $\mathcal{L}_{\text{fermions}}$. Contains the SM terms plus the linearized SMEFT corrections due to field redefinitions and input parameter shifts.
- `LFermionsP`. Same as `LFermions`, but with at least one top quark or W, Z boson replaced with the corresponding dummy field `t1`, `W1`, `Z1`.
- `LYukawa` = $\mathcal{L}_{\text{Yukawa}}$. Contains the SM terms plus the linearized SMEFT corrections due to field redefinitions and input parameter shifts.
- `LYukawaP`. Same as `LYukawa`, but with at least one top quark or Higgs boson replaced with the corresponding dummy field `t1`, `H1`.
- `LSM` = \mathcal{L}_{SM} . The SM Lagrangian without any SMEFT correction.

²⁰All the timings indicated in this section refer to a four-core laptop, with `FeynRules` calculations parallelized.

- **LSMlinear**. The SM lagrangian plus the linearized SMEFT corrections due to field redefinitions and input parameters shifts.
- **LSMloop** = $\mathcal{L}_{\text{SMhloop}}$ as defined in eq. (5.1).
- **LSMloopP**. Same as **LSMloop**, but with at least one Higgs or Z boson replaced with the corresponding dummy field **H1**, **Z1**.
- **LSMincl** = **LSMlinear** + **LSMloop**
- **L6c1[n]** = $\mathcal{L}_6^{(n)}$ with $n = 1 \dots 8$.
 For class 4, the definition is split into **L6c14**, **L6c14cpv** containing only the CP-even and -odd terms respectively.
 For class 8, sub-Lagrangians **L6c18a** ... **L6c18d** are defined in addition.
- **L6no4f** = $\sum_{n=1}^7 \mathcal{L}_6^{(n)}$
- **L6** = $\sum_{n=1}^8 \mathcal{L}_6^{(n)}$
- **LSMEFT** = **LSMincl** + **L6**

The last 3 Lagrangians contain extremely long expressions. It is strongly recommended to use them with care and avoid calling these variables unless strictly necessary.

The parameters notation in the code is provided in appendix D. In addition, the following parameters lists are defined in all models:

- **WC6**. The list of all Wilson coefficients.
 In the **general** model the list **WC6indices** is defined in addition. In this case **WC6** contains e.g. **cHuIm11**, **cHuIm33**, while **WC6indices** contains **cHu[ff1_,ff2_]** with blank flavor indices.
- **shifts**. The list of all shift parameters, such as **dGf**, **dMZ2**, **dgw**, **dg1** etc. The complete list is given in table 17.
- **d6pars**. List of all SMEFT quantities, including Wilson coefficients with and without free indices, and shifts.

Two handy functions are also defined:

- **LinearWC[x_]**. Expands the expression **x** to linear order in the SMEFT parameters (Wilson coefficients and shifts).
- **SMLimit[x_]**. Returns the SM limit of the expression **x**, setting to zero all the **d6pars**.
- **relativeVariation[x_]**. Returns $x / \text{SMLimit}[x]$.
- **SimplifyWC[x_]**. Returns the expression **x** in a form that collects the contributions from each SMEFT parameter (both Wilson coefficients and shifts).

By default, all the Feynman rules are printed out in an input scheme-independent form. The expressions in terms of Wilson coefficients are recovered via a replacement rule that should be applied with `ReplaceRepeated`, as in:

```
In[* ]:= dgw // . MwShifts // WCsimplify
```

$$\text{Out[*]} = -\frac{c_{\text{HL}}^3 \hat{V}^2}{\Lambda^2} + \frac{c_{\text{ll}}^{\text{Prime}} \hat{V}^2}{2 \Lambda^2}$$

if the model was imported with `Scheme = MwScheme`, or

```
In[* ]:= dgw // . alphaShifts // WCsimplify
```

$$\text{Out[*]} = \frac{c_{\text{ll}}^{\text{Prime}} c_{\theta}^2 \hat{V}^2}{2 \Lambda^2 (1 - 2 s_{\theta}^2)} + \frac{c_{\text{HD}} c_{\theta}^2 \hat{V}^2}{4 \Lambda^2 (-1 + 2 s_{\theta}^2)} + \frac{c_{\text{HL}}^3 c_{\theta}^2 \hat{V}^2}{\Lambda^2 (-1 + 2 s_{\theta}^2)} + \frac{c_{\text{HWB}} c_{\theta} s_{\theta} \hat{V}^2}{\Lambda^2 (-1 + 2 s_{\theta}^2)}$$

if the model was imported with `Scheme = alphaScheme`. For consistency, only the appropriate replacement list is defined in each case. All Feynman rules are printed in unitary gauge. It is possible to switch to Feynman gauge by changing the flag `FeynmanGauge` to `True` at any time. However, we caution the user that the SMEFT contributions to the Goldstone and ghost Lagrangians are not fully implemented. In particular, gauge fixing terms have been omitted and the Goldstone kinetic terms are *not* canonically normalized.

The operators' names start with `0` and their definitions carry free flavor indices (eg. `0Hu[ff1_,ff2_]`). In this way they can be shared by multiple setups (the `general`, `U35` and `MFV` models all use a set of definitions, and the `top`, `topU31` models share a separate one). The distinction between flavor assumption is coded through the flavor contractions in the Lagrangian definition, contained in `SMEFTsim_d6_lagrangian.fr`. Therefore it is recommended to isolate each operator through its Wilson coefficient. For instance, the Feynman rules of the operator Q_{Hu} with a $U(3)^5$ flavor symmetry can be printed out via:

```
In[* ]:= Select[L6cl7, !FreeQ[#, cHu] &] // FeynmanRules // Simplify // TableForm
```

Starting Feynman rule calculation.

Expanding the Lagrangian...

Collecting the different structures that enter the vertex.

3 possible non-zero vertices have been found -> starting the computation: 3 / 3.

3 vertices obtained.

```
Out[* ]/TableForm=
```

uq	1	
uq	2	$\frac{i c_{Hu} e \delta_{m_1, m_2} \delta_{f_1, f_2} \gamma^{\mu 5} \cdot P_{+s_1, s_2}}{c_{\theta} \Lambda^2 s_{\theta}}$
H	3	
H	4	
Z	5	
-		
uq	1	
uq	2	$\frac{i c_{Hu} e \hat{v} \delta_{m_1, m_2} \delta_{f_1, f_2} \gamma^{\mu 4} \cdot P_{+s_1, s_2}}{c_{\theta} \Lambda^2 s_{\theta}}$
H	3	
Z	4	
-		
uq	1	
uq	2	$\frac{i c_{Hu} e \hat{v}^2 \delta_{m_1, m_2} \delta_{f_1, f_2} \gamma^{\mu 3} \cdot P_{+s_1, s_2}}{2 c_{\theta} \Lambda^2 s_{\theta}}$
Z	3	

Alternatively, one can reproduce the flavor contraction structure explicitly. In this case:

```
In[*]:= Block[{ff},  $\frac{c_{Hu}}{\Lambda^2 \text{SMEFT}^2}$  OHu[ff, ff] // FeynmanRules // Simplify // TableForm
```

Starting Feynman rule calculation.

Expanding the Lagrangian...

Collecting the different structures that enter the vertex.

3 possible non-zero vertices have been found -> starting the computation: 3 / 3.

3 vertices obtained.

Out[*]//TableForm=

```

-
uq 1
uq 2  $\frac{i c_{Hu} e \delta_{m_1, m_2} \delta_{f_1, f_2} \gamma^{\mu 5} . P_{+s_1, s_2}}{c_{\theta} \Lambda^2 s_{\theta}}$ 
H 3
H 4
Z 5

-
uq 1
uq 2  $\frac{i c_{Hu} e \hat{v} \delta_{m_1, m_2} \delta_{f_1, f_2} \gamma^{\mu 4} . P_{+s_1, s_2}}{c_{\theta} \Lambda^2 s_{\theta}}$ 
H 3
Z 4

-
uq 1
uq 2  $\frac{i c_{Hu} e \hat{v}^2 \delta_{m_1, m_2} \delta_{f_1, f_2} \gamma^{\mu 3} . P_{+s_1, s_2}}{2 c_{\theta} \Lambda^2 s_{\theta}}$ 
Z 3
    
```

or print the result for one flavor entry only:

```
In[*]:=  $\frac{c_{Hu}}{\Lambda^2 \text{SMEFT}^2}$  OHu[1, 1] // FeynmanRules // Simplify // TableForm
```

Starting Feynman rule calculation.

Expanding the Lagrangian...

Collecting the different structures that enter the vertex.

3 possible non-zero vertices have been found -> starting the computation: 3 / 3.

3 vertices obtained.

Out[*]//TableForm=

```

-
u 1
u 2  $\frac{i c_{Hu} e \delta_{m_1, m_2} \gamma^{\mu 5} . P_{+s_1, s_2}}{c_{\theta} \Lambda^2 s_{\theta}}$ 
H 3
H 4
Z 5

-
u 1
u 2  $\frac{i c_{Hu} e \hat{v} \delta_{m_1, m_2} \gamma^{\mu 4} . P_{+s_1, s_2}}{c_{\theta} \Lambda^2 s_{\theta}}$ 
H 3
Z 4

-
u 1
u 2  $\frac{i c_{Hu} e \hat{v}^2 \delta_{m_1, m_2} \gamma^{\mu 3} . P_{+s_1, s_2}}{2 c_{\theta} \Lambda^2 s_{\theta}}$ 
Z 3
    
```

The Feynman rules of the SM Higgs loop Lagrangian are quite complex, especially in the Higgs-gluon operators' case. Their evaluation with the `FeynmanRules` command can be extremely slow and take up to a few hours for the most complex vertices. In order to facilitate their access, the file `SMEFTsim_SMHloop_FRs.nb` is provided, that contains pre-exported expressions. They can be accessed from another notebook via

```
SetDirectory ["PATH/TO/SMEFTSIM/DIR"];
NotebookEvaluate ["SMEFTsim_SMHloop_FRs.nb"];
```

Without producing any output, this will define the objects `lhloop5S`, `lhloop5PS`. The former is a list of all Feynman rules from $\mathcal{L}_{\text{SMhloop}}$, for vertices with up to 5 legs. The latter is the same, but with at least one Higgs boson replaced by the dummy field `H1`. For instance, the first entry is the $h\gamma\gamma$ vertex:

`In[*]:= lhloop5S[[1]]`

$$\text{Out[*]} = \left\{ \{A, 1\}, \{A, 2\}, \{H, 3\}, \frac{4 i g_{H\gamma\gamma} p_1^{\mu_2} p_2^{\mu_1}}{\hat{v}} - \frac{4 i g_{H\gamma\gamma} \eta_{\mu_1, \mu_2} p_1 \cdot p_2}{\hat{v}} \right\}$$

and

`In[*]:= lhloop5PS[[1]]`

$$\text{Out[*]} = \left\{ \{A, 1\}, \{A, 2\}, \{H1, 3\}, \frac{4 i g_{H\gamma\gamma} \Delta_{\text{prop.}} p_1^{\mu_2} p_2^{\mu_1}}{\hat{v}} - \frac{4 i g_{H\gamma\gamma} \Delta_{\text{prop.}} \eta_{\mu_1, \mu_2} p_1 \cdot p_2}{\hat{v}} \right\}$$

is explicitly proportional to the `propCorr` parameter (shown here as $\Delta_{\text{prop.}}$), see section 6.1, 8.3.

Finally, it is *not* recommended to export the UFO models independently, unless only a small subset of operators is included. The UFOs provided in the GitHub repository have been exported in a specific, optimized way and the python files have been manipulated a posteriori in order to introduce the modified form of the propagators and to define the Higgs decay width such that the fermionic decay contributions to $\delta\Gamma_h$ is removed whenever the corresponding Yukawa coupling is set to zero. The notebook with the original export procedure is available upon request.

8 Usage in MadGraph5_aMC@NLO

This section provides recommendations for the use of `SMEFTsim` in `MadGraph5_aMC@NLO`. It is in no way meant as a manual for the functionalities of `MadGraph5_aMC@NLO` itself, for which we defer the reader to the appropriate references, see e.g. [46, 110, 111].

The `SMEFTsim` package provides 10 pre-exported UFO models, one for each flavor setup and input parameter scheme. Each of them contains the full $\mathcal{L}_{\text{SMEFT}}$ defined in section 1 and 3 and $\mathcal{L}_{\text{SMhloop}}$ defined in section 5. The manipulations and redefinitions described in section 2 and 4 have been consistently applied. The vertices contained in the models are derived in unitary gauge and the ghost fields have been removed: `SMEFTsim` is designed for LO event generation and does not support the NLO syntax. A list of the SMEFT parameters defined in the codes is provided in appendix D, with a mapping to the notation used in this notes. All the UFO models have been validated following the recommendations in ref. [47], as detailed in appendix F.

Although the selection of an appropriate model is of course up to the taste of the user, each flavor setup is meant to optimize the parameterization of a certain class of effects in the SMEFT. For instance, the `top` and `topU31` models are designed to single out the couplings of the top and bottom quarks [36], and they only differ in that the `top` case provides more freedom to distinguish the lepton flavors. The `U35` models allow one to work with a minimal number of parameters and are recommended for flavor-blind processes or whenever the flavor structure can be assumed to be strictly SM-like. At the other side of the spectrum, the `general` models provide maximal freedom and can be used to study flavor-violating processes or to realize arbitrary flavor structures beyond those implemented. An operator by operator comparison of the different flavor structures is provided in appendix C.

As discussed in section 6, the use of the $\{m_W, m_Z, G_F\}$ input scheme is particularly recommended for processes involving W bosons, as it avoids the problematic introduction of SMEFT corrections to the W pole mass. The $\{\alpha_{\text{em}}, m_Z, G_F\}$ and $\{m_W, m_Z, G_F\}$ input scheme implementations are expected to give results that differ most significantly in the dependence on the Wilson coefficients C_{HWB} , C_{HD} , $C_{HI}^{(3)}$, C'_u (or $(C_{HI}^{(3)})_{11}$, $(C_{HI}^{(3)})_{22}$, $(C_u)_{1221}$ in models with explicit flavor indices) and in the presence/absence of corrections to either m_W or α_{em} , as described in section 4.2. Further numerical differences affecting both EFT and SM predictions can be present, due to the different definition of the SM parameters in the two cases. These are generally subleading.

8.1 Parameter cards and restrictions

The model parameters are grouped in blocks, that are explicitly shown in the parameter cards. Besides the usual ones (`SMINPUTS`, `MASS`, `DECAY`, `YUKAWA`, `CKMBLOCK`), the parameter card of each `SMEFTsim` UFO model contains the blocks:

`SMEFTCUTOFF` — the parameter Λ . By default this is 1 TeV.

`SMEFT` — the CP-conserving Wilson coefficients, with default value 0.

`SMEFTCPV` — the CP-violating Wilson coefficients, with default value 0.

This block is absent in the MFV models.

`SMEFTFV` — the (ΔC_α) parameters of the MFV setup, with default value 0.

This block is only present in MFV models.

`SWITCHES` — the parameter `linearPropCorrections`, that can be used to switch ON/OFF the linearization of SMEFT corrections in the propagators. The default value is 0 (OFF).

The `SMINPUTS` block contains G_F , α_s and either α_{em} or m_W depending on the input scheme.

The use of restriction cards allows one to reduce the number of diagrams generated for a given process. Two restriction cards are provided by default with each UFO:

- `restrict_massless.dat`. The masses and Yukawa couplings of all fermions, except the top and bottom quarks, are set to 0. The CKM matrix is set to the identity. The Wilson coefficients are set to arbitrary numerical values.
- `restrict_SMLimit_massless.dat`. As in `restrict_massless.dat`, but with all Wilson coefficients set to 0.

The restrictions should be applied at the stage where the model is imported, e.g.:

```
import model SMEFTsim_U35_MwScheme_UFO-massless
```

In this way, all the parameters that are set to either 0 or 1 in the restriction are fixed to their value and cannot be edited further. Sets of parameters that are assigned an identical value in the restriction are fixed to be identical: while their numerical value can still be edited, they cannot be disentangled from one another. Diagrams that are proportional to a vanishing parameter will not be generated.

The use of one of the `massless` restrictions is recommended for LHC studies, because it simplifies significantly the calculations. There are of course several possible strategies for the use of these restrictions in `MadGraph5_aMC@NLO`: for instance, one can create a modified version of `restrict_SMLimit_massless.dat` turning on one Wilson coefficient with some arbitrary value $\neq 0$. Importing the model with this modified restriction allows one to generate events with the chosen coefficient only, while all the other operators are forbidden. Alternatively, if the model is imported with `restrict_massless.dat`, all the Wilson coefficients are retained: all the allowed SMEFT diagrams will be generated and all the parameters can be freely edited at the event generation stage. Note that, to achieve this, all the Wilson coefficients in `restrict_massless.dat` are assigned different non-vanishing and non-unitary values,²¹ that will need to be changed prior to the event generation. To simplify this operation, a “restricted” parameter card `param_card_massless.dat` is provided in the UFO, where all the Wilson coefficients are set to 0. This card can be directly copied in the `PROC/Cards/` directory of the exported process and modified at will.

8.2 Interaction orders

A standard feature of UFO models is that every coupling parameter is assigned an interaction order, i.e. a “flag” that provides control on the number of coupling insertions in generated Feynman diagrams. Each parameter carries an arbitrary number of interaction orders.

²¹In previous versions of `SMEFTsim`, the Wilson coefficients in the restriction cards were all set to the special value `9.999999e-01`, that in principle allows one to set the parameters to 1 without *fixing* their value. However, this syntax is not fully supported by `MadGraph5_aMC@NLO`, and is occasionally source of unexpected numerical behavior in UFO models with a very large number of parameters, such as `SMEFTsim general` or `MFV`.

Order	Parameters assigned
QED = 1	e, g_W, g_1, Y_ψ
QED = 2	$G_F, \hat{\lambda}, \alpha_{\text{em}}$
QED = -1	\hat{v}, Λ
QCD = 1	g_s
QCD = 2	α_s
SMHLOOP = 1	$g_{H\gamma\gamma}, g_{HZ\gamma}, g_{Hgg}^{(1)} \dots g_{Hgg}^{(5)}$
NP = 1	all Wilson coefficients and shifts, except <code>propCorr</code>
NPcpv = 1 (*)	all CP-violating Wilson coefficients
NPfv = 1 (*)	all (ΔC) parameters in MFV models
NPc[a] = 1 (*)	all the CP components and flavor indices of the Wilson coefficient C_α
NPprop = 1 (*)	<code>propCorr</code>

Table 7. Interaction orders defined in the `SMEFTsim` UFO models. Those marked with a (*) are new in version 3.0. The string [a] stands for a generic root name of a Wilson coefficient, as listed in appendix D.

8.2.1 Definitions

In the `SMEFTsim` UFO models the interaction orders are assigned as reported in table 7.

The orders QED and QCD are assigned as customary in the standard SM UFO implementations, with the exceptions of the SMEFT cutoff Λ , that has been assigned QED=-1 such that the combination (\hat{v}/Λ) is order-less, and of the Wilson coefficient C_H , that has been assigned QED=1. This prevents the C_H correction to the h^3 interaction, that is proportional to $\hat{v}^3 C_H / \Lambda^2$ (see eq. (4.44)), from having overall order QED = -1.

The interaction order SMHLOOP labels the SM loop-generated Higgs interactions introduced in section 5. Since by definition they are proportional to the SM gauge couplings, the $g_{Hgg}^{(k)}$ parameters additionally carry QCD=2 and the $g_{H\gamma\gamma}, g_{HZ\gamma}$ parameters carry QED=2.

The interaction order NP (New Physics) is assigned to all the Wilson coefficients and shifts indistinctly. In addition, starting from version 3.0, individual interaction orders have been introduced for each effective operators. The same order NPc[a] is assigned to all the associated CP-conserving and violating parameters, irrespective of the flavor indices carried. For instance, in the `top` models, the parameters $\text{Re}(C_{eH})_{pp}, \text{Im}(C_{eH})_{pp}$ for $p = \{1, 2, 3\}$ all have order NPceH=1. In the `U35`, `MFV`, `top` and `topU31` models, distinct interaction orders are assigned to independent flavor contractions. For instance C_{ll} and C'_{ll} have orders NPc11 and NPc111 respectively. In the `top` models, the parameters $(C_{ll})_{pprr}$ have order NPc11, while the $(C_{ll})_{pprrp}$ contractions have order NPc111, etc. In most cases the label [a] coincides with the name root of the associated Wilson coefficient, that can be read off from the tables in appendix D. If in doubt, the user can resort to the `.fr` source files or check explicitly the `couplings.py` file to identify the exact orders assigned to a given parameter or coupling.

All the CP-violating parameters, that belong to the SMEFTCPV block, have an order $\text{NPcpv}=1$. Analogously, all the (ΔC_α) quantities in the MFV models, that belong to the SMEFTFV block, have an order $\text{NPfv}=1$.

Finally, the order NPprop labels the interactions of the dummy fields W', Z', h', t' carrying linearized propagator corrections, see section 6.1. It is carried by a dummy internal parameter propCorr that only takes values 0/1, when the `linearPropCorrections` switch is set to 0/a nonzero value. Its application is discussed in the next subsection. By default, the interaction order NPprop is “switched off”, as it is assigned an upper limit of 0 interactions, that can be lifted as shown below. No upper limit is set for the other orders.

The interaction orders SMHLOOP , NP , NPprop , NPcpv , NPfv have been assigned hierarchy 99. `MadGraph5_aMC@NLO` will therefore generally avoid insertions of the associated vertices, unless these orders are specified.

8.2.2 Recommended use

Interaction orders are specified at the stage of process generation in `MadGraph5_aMC@NLO`, e.g.:

```
generate p p > mu+ mu- SMHLOOP=0 NPprop=0 NP=1 NP^2==1
```

where `=` is equivalent to `<=`, while `==` selects uniquely the order specified. The syntax `XX=n` acts at the amplitude level, i.e. it specifies the total number of couplings with order `XX` to be inserted in each Feynman diagram. The syntax `XX^2` acts instead at the *squared* amplitude level. This functionality works very nicely for EFT studies, as it allows one to disentangle contributions at different orders in the expansion.

Although *a priori* `SMEFTsim` can be used for computations to any allowed order in Λ , it implements the SMEFT Lagrangian consistently expanded only up to $\mathcal{O}(\Lambda^{-2})$. This means that any `SMEFTsim` prediction beyond this order is necessarily incomplete in the Effective Theory. It is worth noting that this statement does not concern only higher dimensional operators in $\mathcal{L}_{\text{SMEFT}}$, but also affects the dependence on some of the Wilson coefficients in \mathcal{L}_6 . For instance, it was stressed at multiple stages in sections 2 and 4 that terms of order Λ^{-4} or higher were neglected in the field and parameter redefinitions performed, as well as in the treatment of input parameters. The impact of these \mathcal{L}_6 contributions has been discussed in ref. [112] for the case of $\mathcal{O}(\Lambda^{-4})$ corrections to $1 \rightarrow 2$ decays, using the `geoSMEFT` formalism [62].

Complete results truncated at $\mathcal{O}(\Lambda^{-2})$ can be obtained with the syntax `NP<=1 NP^2<=1`, that retains only SM plus SM- \mathcal{L}_6 interference contributions. Contributions of order Λ^{-4} stemming from the square of an $\mathcal{O}(\Lambda^{-2})$ amplitude, although incomplete, are also commonly included in the SMEFT calculations: they are selected with `NP=1 NP^2==2` or just `NP==1`. For the reasons above, it is generally recommended to use the specification `NP<=1` (or `NP=1`) for any process, to limit the number of EFT insertions to one per Feynman diagram. A generic observable computed in this way will have the form

$$\sigma = \sigma_{\text{SM}} + \sum_{\alpha} \sigma_{\alpha} \bar{C}_{\alpha} + \sum_{\alpha, \beta} \sigma_{\alpha\beta} \bar{C}_{\alpha} \bar{C}_{\beta}, \tag{8.1}$$

	σ_{SM}	σ_{α}	σ_{β}	$\sigma_{\alpha\alpha}$	$\sigma_{\beta\beta}$	$\sigma_{\alpha\beta}$
NP=0	✓					
NP<=1	✓	✓	✓	✓	✓	✓
NP==1				✓	✓	✓
NP<=1 NP^2<=1	✓	✓	✓			
NP<=1 NP^2==1		✓	✓			
NP<=1 NPc [a]^2<=1	✓	✓				✓
NP<=1 NPc [a]^2<=1 NPc [b]^2<=1	✓	✓	✓			✓
NP<=1 NPc [a]==1		✓		✓		
NP<=1 NPc [a]^2==1		✓				✓
NP<=1 NPc [a]^2==2				✓		
NP<=1 NP^2==1 NPc [a]^2==1		✓				
NP<=1 NP^2==2 NPc [a]^2==1						✓

Table 8. Examples of interaction-order syntax that select different EFT contributions to a generic observable with the dependence given in eq. (8.1), for the case of 2 Wilson coefficients C_{α}, C_{β} .

where $\sigma_{\text{SM}}, \sigma_{\alpha}, \sigma_{\alpha\beta}$ intuitively denote the SM, interference and quadratic contributions respectively. σ_{SM} and $\sigma_{\alpha\alpha}$ are always positive quantities, while $\sigma_{\alpha}, \sigma_{\alpha\beta}$ with $\alpha \neq \beta$ can take negative values. Table 8 shows examples of how the interaction order syntax can be used to disentangle these contributions, for a simple case with two Wilson coefficients. The expressions directly generalize if three or more parameters are present.

Finally, as discussed in section 5, the loop-generated SM Higgs couplings implemented in SMEFTsim are defined in the $m_t \rightarrow \infty$ limit, and their use should be limited to (on-shell) Higgs production and decay processes. Outside of this regime, it is strongly recommended to use SMHLOOP=0.

8.3 Propagator corrections and decay widths

As discussed in section 6, SMEFT corrections are generally present in the propagators of unstable particles, due to $d = 6$ operators modifying their masses and/or decay widths. Section 6.1 outlined two alternative methods for estimating these contributions in a given process. In the following we illustrate how they can be implemented in MadGraph5_aMC@NLO.

8.3.1 Method (a): linearized corrections

SMEFTsim 3.0 offers the possibility to linearize propagator corrections as in section 6 for the Z, W, h bosons and for the top quark. The implementation relies on the introduction of dummy fields Z', W', h', t' whose couplings carry interaction order NPprop, which allows one to single them out. For instance, the syntax

```
generate p p > e+ e- SMHLOOP=0 NP==1 NPprop=0
```


selects all diagrams with one effective operator in a *vertex* (including 4-point $\bar{q}qe^+e^-$ interactions) but none in propagators, while

```
generate p p > e+ e- SMHLOOP=0 NP=0 NPprop<=2 NPprop^2==2
generate p p > e+ e- SMHLOOP=0 NP=0 NPprop==2
```

extract the pure SM- \mathcal{L}_6 interference ($\propto \delta\Gamma_Z$) and the quadratic ($\propto (\delta\Gamma_Z)^2$) contributions respectively from corrections to the Z propagator,²² and exclude EFT insertions in vertices. Due to the absence of additional interaction orders, there is unfortunately no equivalent to table 8 in this case: propagator corrections from different operators and EFT orders cannot be disentangled at this level.

Important: in order to avoid unwanted insertions of the dummy fields in standard process generations, the functionality described here has to be activated in 2 steps: (i) in the file `coupling_order.py`, the `expansion_order` option for the order `NPprop` has to be set to a number ≥ 2 (recommended: 99). The default is 0, which forbids dummy interactions completely. (ii) The parameter `linearPropCorrections` in the `param_card` has to be set to a non-zero value. If this is not the case, dummy vertices will be included in the diagrams, but they will be idle, as they are proportional to `propCorr = 0/1` for `linearPropCorrections = 0/non-zero`.

8.3.2 Method (b): full corrections

As an alternative to linearization, propagator corrections can be estimated following more canonical procedures. This generally means computing processes with the propagator forms in eqs. (6.1)–(6.3), with mass and decay parameters that depend explicitly on the Wilson coefficients, to either linear or quadratic order. The resulting process will thus exhibit a non-polynomial dependence on the SMEFT parameters.

The most relevant caveat here is that the implementation of the Wilson coefficient dependence is necessarily different for masses and widths. In the former case, it is possible to define mass parameters as internal and assign them an analytic expression, e.g. $M_W = M_{Wsm} + dMW$, with dMW defined as in (4.33). In `SMEFTsim` the dMW term is only included when `linearPropCorrections = 0`, and switched off otherwise. Note also that the expression of a generic δm is extracted at the Lagrangian level and is purely of $\mathcal{O}(\Lambda^{-2})$.

On the other hand, due to how Monte Carlo generators and their interface to parton shower or decay modules are structured, decay widths cannot be defined as internal parameters in `UFO` models. Therefore the only way their SMEFT expressions can be inserted in the calculation is by letting `MadGraph5_aMC@NLO` compute them, by setting the relevant widths to `Auto` in the `param_card` [106]. The on-the-fly calculation will include all allowed 2-body decays as well as higher multiplicity decays estimated to be numerically relevant, and it will rely on the pre-computed decay results collected in the file `decays.py`, which

²²Remember that `NPprop` counts the number of dummy *vertices*, so, in this case, the order specified is twice the number of dummy propagators.

include $\mathcal{O}(\Lambda^{-4})$ terms.²³ Note that, with this procedure, the decay widths will need to be re-evaluated every time the value of a relevant Wilson coefficient is modified.

With both mass and width corrections evaluated as above, the denominator of a generic propagator has the form

$$q^2 - (m^{\text{SM}} + \delta m)^2 + i(m^{\text{SM}} + \delta m)(\Gamma^{\text{SM}} + \delta^{(1)}\Gamma + \delta^{(2)}\Gamma), \quad (8.2)$$

where δm and $\delta^{(1)}\Gamma$ are of $\mathcal{O}(\Lambda^{-2})$ and $\delta^{(2)}\Gamma$ is of $\mathcal{O}(\Lambda^{-4})$. Eq. (8.2) contains therefore terms up to $\mathcal{O}(\Lambda^{-6})$. An observable computed for a process with k internal lines corrected in this way, will contain terms up to $\mathcal{O}(\Lambda^{-12k})$ at the denominator.

In principle this functional dependence can be reconstructed fitting the appropriate rational function to a sufficient number of benchmark points. However, it is recommended to reduce the proliferation of higher-order terms in the propagators, by evaluating mass and width corrections separately and by treating propagator corrections to different internal states individually, whenever possible. This is achieved avoiding to switch on at the same time the `linearPropCorrections` flag (that turns on δm_W) and the `Auto` computation of a decay width, or of two decay widths simultaneously.

8.4 Example: Higgs production and decay including W, Z propagator corrections

As a practical example for the use of `SMEFTsim` in `MadGraph5_aMC@NLO`, we compute SMEFT corrections to Higgs production and decay processes that are mediated by W, Z exchange, as an illustration of the propagator corrections feature.

8.4.1 STXS for $\bar{q}q \rightarrow h\bar{q}q$

We consider two bins of the stage 1.1 Simplified Template Cross Section (STXS) parameterization [113–116], for the EW $\bar{q}q \rightarrow h\bar{q}q$ production channel at low Higgs p_T .

They are defined by the cuts [115]:

$$\begin{array}{l|l} \text{VBF-like} & 350 \text{ GeV} < m_{jj} & p_T(h) < 200 \text{ GeV}, |y_h| < 2.5 \\ \text{VH-like} & 60 \text{ GeV} < m_{jj} < 120 \text{ GeV} & p_T(h) < 200 \text{ GeV}, |y_h| < 2.5. \end{array}$$

with y_h the rapidity of the Higgs boson. In each bin, the Higgs production cross section in the SMEFT can be parameterized as:

$$\sigma_{\text{SMEFT}} = \sigma_{\text{SM}} + \sum_{\alpha} \sigma_{\alpha} \bar{C}_{\alpha} + \mathcal{O}(\Lambda^{-4}). \quad (8.3)$$

Table 9 reports the values of $\sigma_{\alpha}/\sigma_{\text{SM}}$ for the relevant fermionic operators, computed at parton level using `SMEFTsim` in the `U35` flavor-symmetric, $\{m_W, m_Z, G_F\}$ input scheme version. The following procedure was followed:

1. 50000 events are generated for each bin in `MadGraph5_aMC@NLO`, for $\sqrt{s} = 13 \text{ TeV}$. The syntax used is

²³Since only $1 \rightarrow 2$ decays are included, these results consistently stem from the square of $\mathcal{O}(\Lambda^{-2})$ amplitudes. Dummy fields are not included in the pre-computed decay widths.

```
import model SMEFTsim_U35_MwScheme_UF0-vbf
generate q q > h q q QCD=0 NP=0 SMHLOOP=0 NPprop=0
```

where the `-vbf` flag indicates that the model is imported with a custom restriction card `restrict_vbf.dat`, that in this case sets to zero the masses and Yukawa couplings of all fermions except the bottom and top quarks, as well as all the Wilson coefficients that are known not to contribute to the process. The remaining ones are set to a random non-zero value in this card. The STXS defining cuts in $p_T(h)$ and y_h cuts are implemented at the level of the `run_card.dat` in `MadGraph5_aMC@NLO`, while the invariant mass cuts are applied when analyzing the events a posteriori. This gives the tree-level SM cross sections

$$\sigma_{SM}^{VBF} = 1.56 \text{ pb}, \quad \sigma_{SM}^{VH} = 0.67 \text{ pb}. \tag{8.4}$$

2. The events are reweighted using the reweight module in `MadGraph5_aMC@NLO` [117]. Individual weights are computed for each Wilson coefficient, splitting contributions from operator insertions in the vertices (labeled as “direct”) and from insertions in the W, Z propagators (labeled as “propagator”). This is done setting each coefficient to 1 and the SMEFT cutoff scale `LambdaSMEFT` to 1 TeV. For instance, for the C'_u parameter, the `reweight_card.dat` for the direct contributions is

```
change process q q > h q q QCD=0 NP=1 NP^2==1 NPprop=0 SMHLOOP=0
launch --rwgt_name=SMEFTsim-cl11-direct
set cll1 1
set cH11 0
set cH13 0
set cHe 0
set cHq1 0
set cHq3 0
set cHu 0
set cHd 0
done
```

For estimating the pure propagator contributions the first two lines are replaced with

```
change process q q > h q q QCD=0 NP=0 NPprop=2 NPprop^2==2 SMHLOOP=0
launch --rwgt_name=SMEFTsim-cl11-propagator
```

Analyzing the reweighted events gives $(\sigma_\alpha \frac{\hat{v}^2}{\Lambda^2})$ for each C_α . The numbers in table 9 are finally obtained dividing by σ_{SM} and normalizing to \bar{C}_α .

The results show that propagator corrections are negligible in the VBF regime, where the relative SMEFT corrections to the cross section is

$$\frac{\sigma_{SMEFT}}{\sigma_{SM}} = 1 - 3 \cdot 10^{-4} \frac{\delta\Gamma_Z}{\Gamma_Z^{SM}} - 7 \cdot 10^{-4} \frac{\delta\Gamma_W}{\Gamma_W^{SM}} + \text{direct}. \tag{8.5}$$

	$\bar{q}q \rightarrow h\bar{q}q$ VBF-like		$\bar{q}q \rightarrow h\bar{q}q$ VH-like		$h \rightarrow e^+e^-\mu^+\mu^-$	
	direct	propagators	direct	propagators	direct	propagators
\bar{C}_{He}		$5.32 \cdot 10^{-5}$		0.0526	-1.724	0.153
$\bar{C}_{Hl}^{(1)}$		$5.32 \cdot 10^{-5}$		0.0526	2.144	0.153
$\bar{C}_{Hl}^{(3)}$	-6	$1.351 \cdot 10^{-3}$	-6	1.258	-3.856	1.147
$\bar{C}_{Hq}^{(1)}$	0.109	$-1.363 \cdot 10^{-4}$	-0.197	-0.135		-0.39
$\bar{C}_{Hq}^{(3)}$	-5.345	$-1.423 \cdot 10^{-3}$	25.66	-1.329		-1.353
\bar{C}_{Hu}	-0.323	$-7.092 \cdot 10^{-5}$	1.926	-0.070		-0.203
\bar{C}_{Hd}	0.103	$5.24 \cdot 10^{-5}$	-0.608	0.0518		0.150
\bar{C}'_u	3	$-1. \cdot 10^{-3}$	3	-0.936	3	-0.839

Table 9. Values of σ_a/σ_{SM} for the relevant fermionic Wilson coefficients \bar{C}_α contributing to $\bar{q}q \rightarrow h\bar{q}q$ and $h \rightarrow e^+e^-\mu^+\mu^-$. For the first two columns σ is the cross section in the VBF-like and VH-like STXS bins defined in the text, while in the third the numbers refer to the partial decay width. The results are given for $\{m_W, m_Z, G_F\}$ inputs with a $U(3)^5$ flavor symmetry, and neglecting all fermion masses. The “direct” contributions stem from operator insertions in vertices, including parameter shifts, while the “propagator” ones stem from the corrections to the W, Z decay width in internal propagators. The lines highlighted in color are those for which the latter are most relevant.

This is expected, as the intermediate bosons in t -channel are mostly off-shell in this process. In the VH bin, on the other hand, one vector boson can be on-shell, which enhances the propagator effects. In this case, the relative SMEFT correction is

$$\frac{\sigma_{SMEFT}}{\sigma_{SM}} = 1 - 0.29 \frac{\delta\Gamma_Z}{\Gamma_Z^{SM}} - 0.65 \frac{\delta\Gamma_W}{\Gamma_W^{SM}} + \text{direct}, \tag{8.6}$$

where the numerical prefactors reflect the proportions of W and Z bosons produced. In fact, the largest numerical effects in table 9 are observed in the operators entering $\delta\Gamma_W$.

8.4.2 $h \rightarrow e^+e^-\mu^+\mu^-$

An analysis of the Z -mediated Higgs decay $h \rightarrow e^+e^-\mu^+\mu^-$ was performed following a procedure analogous to the one described for $\bar{q}q \rightarrow h\bar{q}q$. In the decay case, one Z boson is always on-shell, leading to significant contributions from the intermediate Z propagator. The relative SMEFT correction to the decay width is found to be

$$\frac{\Gamma_{SMEFT}}{\Gamma_{SM}} = 1 - 0.84 \frac{\delta\Gamma_Z}{\Gamma_Z^{SM}} + \text{direct}. \tag{8.7}$$

The breakdown into fermionic Wilson coefficients is given in table 9 and it agrees with the analytic results of ref. [57].

9 Summary

The `SMEFTsim` package contains models in `FeynRules` and in the `UFO` format, that implement the complete Warsaw basis of dimension six operators, under different flavor assumptions and with different choices of the input quantities for the EW sector. Its main scope is the Monte Carlo simulation of LHC processes in the SMEFT, but it can also be employed for simple analytic calculations, exploiting the `FeynRules` interface in `Mathematica`.

This work reviewed the theoretical elements that are implemented in `SMEFTsim` and presented the improvements in version 3.0. The most significant changes compared to previous releases are the addition of two new flavor structures for top quark physics, the implementation of a brand new tool for the inclusion of SMEFT corrections in the propagator of unstable particles and the general improvement of the code, particularly of the parameterization of Higgs-gluon interactions in the SM.

As in previous versions, `SMEFTsim` 3.0 supports the `WCxf` exchange format [22]. The corresponding interface will be updated shortly after the code release. Finally, support for the translation of the `UFO` models to `python3` will be provided in the near future.

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A Analytic expressions of decay width corrections

This appendix reports analytic expressions for the decay widths of the W, Z, h, t particles in the SMEFT, that are implemented in the tool documented in section 6.1.

Only linear terms in the Wilson coefficients are retained. CKM mixing and all fermion masses except m_b and m_t are neglected, unless otherwise specified. Results are reported in the flavor-general setup, and they can be mapped to the symmetric scenarios using the tables in appendix C. We use an input-scheme independent notation. Scheme-specific results can be obtained replacing the generic shifts $\delta g_W, \delta g_1, \delta m_W^2$ with the expressions reported in section 4.2.1 or 4.2.2. The quantities $\Delta G_F, \Delta m_Z^2, \Delta \kappa_H$ are defined in table 4. Finally, we use the \bar{C} notation defined in eq. (2.4) and the hat notation defined in section 4.

A.1 Z boson

The Z boson couplings to a fermion pair $\psi\bar{\psi}$ are

$$g_{\psi L} = T_3^\psi - Q_\psi s_\theta^2, \quad g_{\psi R} = -Q_\psi s_\theta^2, \quad (\text{A.1})$$

where $T_3^\psi = \pm 1/2$ is the isospin eigenvalue and Q_ψ is the electric charge of the fermion ψ . We also define:

$$\Delta_{\nu,pp}^L = (\bar{C}_{Hl}^{(1)})_{pp} - (\bar{C}_{Hl}^{(3)})_{pp}, \quad (\text{A.2})$$

$$\Delta_{e,pp}^L = (\bar{C}_{Hl}^{(1)})_{pp} + (\bar{C}_{Hl}^{(3)})_{pp}, \quad \Delta_{e,pp}^R = (\bar{C}_{He})_{pp}, \quad (\text{A.3})$$

$$\Delta_{u,pp}^L = (\bar{C}_{Hq}^{(1)})_{pp} - (\bar{C}_{Hq}^{(3)})_{pp}, \quad \Delta_{u,pp}^R = (\bar{C}_{Hu})_{pp}, \quad (\text{A.4})$$

$$\Delta_{d,pp}^L = (\bar{C}_{Hq}^{(1)})_{pp} + (\bar{C}_{Hq}^{(3)})_{pp}, \quad \Delta_{d,pp}^R = (\bar{C}_{Hd})_{pp}. \quad (\text{A.5})$$

At tree level in the SM, the partial decay width of the Z boson into a $\bar{\psi}_p\psi_p$ pair, with $\psi = \{\nu, l^-, u, d\}$ and flavor p , is

$$\Gamma_{Z \rightarrow \bar{\psi}_p\psi_p}^{\text{SM}} = \frac{G_F m_Z^3}{3\sqrt{2}\pi} N_C^\psi [g_{\psi L}^2 + g_{\psi R}^2], \quad (\text{A.6})$$

where N_C^ψ is the number of colors of the fermion species ψ .

The relative SMEFT correction to a partial width can be inferred differentiating in the $g_{\psi L}, g_{\psi R}$ couplings and inserting the expressions of their SMEFT shifts:

$$\delta g_{\psi L} = g_{\psi L} \delta g_Z - Q_\psi \delta s_\theta^2 - \frac{\Delta_{\psi,pp}^L}{2}, \quad (\text{A.7})$$

$$\delta g_{\psi R} = g_{\psi R} \delta g_Z - Q_\psi \delta s_\theta^2 - \frac{\Delta_{\psi,pp}^R}{2}, \quad (\text{A.8})$$

with

$$\delta g_Z = \frac{\delta g_W}{\hat{g}_W} + \frac{\delta s_\theta^2}{2c_\theta^2} + \frac{s_\theta}{2c_\theta} \bar{C}_{HWB}. \quad (\text{A.9})$$

Using the expression of δs_θ^2 provided in eq. (4.32), one obtains

$$\begin{aligned} \frac{\delta \Gamma_{Z \rightarrow \bar{\psi}_p\psi_p}}{\Gamma_{Z \rightarrow \bar{\psi}_p\psi_p}^{\text{SM}}} &= 2c_\theta^2 \left[1 + 2s_\theta^2 Q_\psi \frac{g_{\psi L} + g_{\psi R}}{g_{\psi L}^2 + g_{\psi R}^2} \right] \frac{\delta g_W}{\hat{g}_W} + 2s_\theta^2 \left[1 - 2c_\theta^2 Q_\psi \frac{g_{\psi L} + g_{\psi R}}{g_{\psi L}^2 + g_{\psi R}^2} \right] \frac{\delta g_1}{\hat{g}_1} \\ &+ s_{2\hat{\theta}} \left[1 - c_{2\hat{\theta}} Q_\psi \frac{g_{\psi L} + g_{\psi R}}{g_{\psi L}^2 + g_{\psi R}^2} \right] \bar{C}_{HWB} - \frac{g_{\psi L} \Delta_{\psi,pp}^L + g_{\psi R} \Delta_{\psi,pp}^R}{g_{\psi L}^2 + g_{\psi R}^2}. \end{aligned} \quad (\text{A.10})$$

Flavor violating decays are absent at $\mathcal{O}(\Lambda^{-2})$. As $m_b \neq 0$ is retained, the $Z \rightarrow \bar{b}b$ result contains additional terms. The partial width expression in the SM is

$$\Gamma_{Z \rightarrow \bar{b}b}^{\text{SM}} = \frac{G_F m_Z^3}{\sqrt{2}\pi} [g_{dL}^2 + g_{dR}^2] \sqrt{1 - 4x_b^2} \left[1 - x_b^2 \left(1 - \frac{6 g_{dL} g_{dR}}{g_{dL}^2 + g_{dR}^2} \right) \right], \quad (\text{A.11})$$

with $x_b = m_b/m_Z$. The relative SMEFT correction is

$$\begin{aligned} \frac{\delta\Gamma_{Z\rightarrow\bar{b}b}}{\Gamma_{Z\rightarrow\bar{b}b}^{\text{SM}}} &= \frac{\delta\Gamma_{Z\rightarrow\bar{b}b}}{\Gamma_{Z\rightarrow\bar{b}b}^{\text{SM}}}\Big|_{m_b=0} + \frac{3x_b^2}{(g_{dL}^2 + g_{dR}^2)(1-x_b^2) + 6x_b^2 g_{dL} g_{dR}} \frac{g_{dL}^2 - g_{dR}^2}{g_{dL}^2 + g_{dR}^2} \times \\ &\times \left[-\frac{1}{3} s_{2\hat{\theta}}^2 (g_{dL} - g_{dR}) \left(\frac{\delta g_W}{\hat{g}_W} - \frac{\delta g_1}{\hat{g}_1} - \frac{c_{2\hat{\theta}}}{s_{2\hat{\theta}}} \bar{C}_{HWB} \right) + g_{dL} \Delta_{d,33}^R - g_{dR} \Delta_{d,33}^L \right] \\ &- \frac{6\sqrt{2} x_b (g_{dL}^2 - g_{dR}^2)}{(g_{dL}^2 + g_{dR}^2)(1-x_b^2) + 6x_b^2 g_{dL} g_{dR}} \left[c_{\hat{\theta}} (\bar{C}_{dW})_{33} + s_{\hat{\theta}} (\bar{C}_{dB})_{33} \right], \end{aligned} \quad (\text{A.12})$$

where the first term stands for the contributions in eq. (A.10). The relative SMEFT correction to the total decay width is finally obtained as

$$\frac{\delta\Gamma_Z}{\Gamma_Z^{\text{SM}}} = \sum_f \left[\frac{\delta\Gamma_{Z\rightarrow f}}{\Gamma_{Z\rightarrow f}^{\text{SM}}} \right] \text{Br}_{Z\rightarrow f}^{\text{SM}}, \quad (\text{A.13})$$

with f running over all the allowed fermion pairs and $\text{Br}_{Z\rightarrow f}^{\text{SM}}, \Gamma_Z^{\text{SM}}$ computed directly from the tree level expressions.

A.2 W boson

At tree level in the SM, the partial decay width of the W^+ boson into a fermion pair $f^+ = \{e_p^+ \nu_p, \bar{d}_p u_p\}$ with flavor p is

$$\Gamma_{W^+\rightarrow f^+}^{\text{SM}} = \frac{G_F m_W^3}{6\sqrt{2}\pi} N_C^f, \quad (\text{A.14})$$

with $N_C^f = \{1, 3\}$ the number of colors. Only decays into same-generation fermions are considered here, as CKM mixing is neglected. The relative SMEFT correction for each channel is

$$\frac{\delta\Gamma_{W^+\rightarrow l_p^+ \nu_p}}{\Gamma_{W^+\rightarrow l_p^+ \nu_p}^{\text{SM}}} = 2 \frac{\delta g_W}{\hat{g}_W} + \frac{\delta m_W}{\hat{m}_W} + 2(\bar{C}_{Hl}^{(3)})_{pp}, \quad (\text{A.15})$$

$$\frac{\delta\Gamma_{W^+\rightarrow u_p \bar{d}_p}}{\Gamma_{W^+\rightarrow u_p \bar{d}_p}^{\text{SM}}} = 2 \frac{\delta g_W}{\hat{g}_W} + \frac{\delta m_W}{\hat{m}_W} + 2(\bar{C}_{Hq}^{(3)})_{pp}. \quad (\text{A.16})$$

The total W^+ decay width in the SM is

$$\Gamma_W = \sum_{p=1}^3 \Gamma_{W^+\rightarrow l_p^+ \nu_p} + \sum_{p=1}^2 \Gamma_{W^+\rightarrow u_p \bar{d}_p}. \quad (\text{A.17})$$

Since in this case the branching ratios are simple rational numbers, the relative SMEFT correction simplifies into

$$\frac{\delta\Gamma_W}{\Gamma_W^{\text{SM}}} = \frac{1}{9} \sum_{p=1}^3 \frac{\delta\Gamma_{W^+\rightarrow l_p^+ \nu_p}}{\Gamma_{W^+\rightarrow l_p^+ \nu_p}^{\text{SM}}} + \frac{1}{3} \sum_{p=1}^2 \frac{\delta\Gamma_{W^+\rightarrow \bar{d}_p u_p}}{\Gamma_{W^+\rightarrow \bar{d}_p u_p}^{\text{SM}}} \quad (\text{A.18})$$

$$= 2 \frac{\delta g_W}{\hat{g}_W} + \frac{\delta m_W}{\hat{m}_W} + \frac{2}{9} \sum_{p=1}^3 (\bar{C}_{Hl}^{(3)})_{pp} + \frac{2}{3} \sum_{p=1}^2 (\bar{C}_{Hq}^{(3)})_{pp}. \quad (\text{A.19})$$

A.3 Higgs boson

The SM partial widths for two-body Higgs decays are:

$$\Gamma_{h \rightarrow \bar{\psi}_p \psi_p}^{\text{SM}} = \frac{m_h y_{\psi_p}^2}{16\pi} N_C^\psi [1 - 4x_{\psi_p}]^{3/2}, \quad (\text{A.20})$$

$$\Gamma_{h \rightarrow \gamma\gamma}^{\text{SM}} = \frac{m_h^3}{4\pi\hat{v}^2} g_{H\gamma\gamma}^2, \quad (\text{A.21})$$

$$\Gamma_{h \rightarrow Z\gamma}^{\text{SM}} = \frac{m_h^3}{8\pi\hat{v}^2} g_{HZ\gamma}^2 [1 - x_Z^2]^3, \quad (\text{A.22})$$

$$\Gamma_{h \rightarrow gg}^{\text{SM}} = \frac{m_h^3}{2\pi\hat{v}^2} \left[2g_{Hgg}^{(1)} - \frac{1}{x_t^2} g_{Hgg}^{(2)} \right]^2, \quad (\text{A.23})$$

where $x_{\psi_p} = m_{\psi_p}/m_h$ for a generic fermion ψ_p , $x_Z = m_Z/m_h$, $x_t = m_t/m_h$ and $y_{\psi_p} \equiv (Y_\psi)_{pp}$ is the relevant Yukawa coupling. The relative SMEFT corrections are

$$\frac{\delta\Gamma_{h \rightarrow \bar{\psi}_p \psi_p}}{\Gamma_{h \rightarrow \bar{\psi}_p \psi_p}^{\text{SM}}} = 1 - \Delta G_F + 2\Delta\kappa_H - 2\frac{(\bar{C}_{\psi H})_{pp}}{y_{\psi_p}}, \quad (\text{A.24})$$

$$\frac{\delta\Gamma_{h \rightarrow \gamma\gamma}}{\Gamma_{h \rightarrow \gamma\gamma}^{\text{SM}}} = \frac{2}{g_{H\gamma\gamma}^2} \left[c_\theta^2 \bar{C}_{HB} + s_\theta^2 \bar{C}_{HW} - s_\theta c_\theta \bar{C}_{HWB} \right], \quad (\text{A.25})$$

$$\frac{\delta\Gamma_{h \rightarrow Z\gamma}}{\Gamma_{h \rightarrow Z\gamma}^{\text{SM}}} = \frac{2}{g_{HZ\gamma}^2} \left[s_{2\theta} (\bar{C}_{HW} - \bar{C}_{HB}) - c_{2\theta} \bar{C}_{HWB} \right], \quad (\text{A.26})$$

$$\frac{\delta\Gamma_{h \rightarrow gg}}{\Gamma_{h \rightarrow gg}^{\text{SM}}} = \frac{4\bar{C}_{HG}}{2g_{Hgg}^{(1)} - g_{Hgg}^{(2)}/x_t^2}, \quad (\text{A.27})$$

with the $h\gamma\gamma, hZ\gamma, hgg$ couplings defined in section 5. The `SMEFTsim` implementation retains the masses of the tau lepton, charm and bottom quarks in eq. (A.24). Four-body decays were included neglecting neutral current contributions, CKM mixing and all fermion masses. The analytic expressions were taken from ref. [57] and generalized to all flavor setups.

For each individual decay channel, the partial decay width in the SM is

$$\Gamma_{h \rightarrow f_+ f_-}^{\text{SM}} = N_C^{f_+} N_C^{f_-} \frac{m_W^8 G_F^3}{m_h} \left(4.65 \times 10^{-4} \right), \quad (\text{A.28})$$

with $f_+ = \{l_p^+ \nu_p, \bar{d}_p u_p\}$, $f_- = \{\bar{\nu}_p l_p^-, \bar{u}_p d_p\}$ and $N_C^{f_+}, N_C^{f_-}$ the appropriate color multiplicities. The numerical factor comes from the phase space integration, that is performed taking $\hat{m}_W = 80.387 \text{ GeV}$.

Summing over all allowed flavor combinations, the relative SMEFT corrections are:

$$\frac{\delta\Gamma_{h \rightarrow l^+ \nu \bar{\nu} l^-}}{\Gamma_{h \rightarrow l^+ \nu \bar{\nu} l^-}^{\text{SM}}} = F_1 + \left(\frac{4}{3} - 0.588 \right) \sum_{p=1}^3 (\bar{C}_{Hl}^{(3)})_{pp}, \quad (\text{A.29})$$

$$\frac{\delta\Gamma_{h \rightarrow \bar{u} d \bar{d} u}}{\Gamma_{h \rightarrow \bar{u} d \bar{d} u}^{\text{SM}}} = F_1 + (2 - 0.882) \sum_{p=1}^2 (\bar{C}_{Hq}^{(3)})_{pp}, \quad (\text{A.30})$$

$$\frac{\delta\Gamma_{h \rightarrow l^+ \nu \bar{u} d + \text{h.c.}}}{\Gamma_{h \rightarrow l^+ \nu \bar{u} d + \text{h.c.}}^{\text{SM}}} = F_1 + \left(\frac{2}{3} - 0.294 \right) \sum_{p=1}^3 (\bar{C}_{Hl}^{(3)})_{pp} + (1 - 0.441) \sum_{p=1}^2 (\bar{C}_{Hq}^{(3)})_{pp}, \quad (\text{A.31})$$

with

$$F_1 = -\Delta G_F + 2\Delta\kappa_H + 4\frac{\delta g_W}{\hat{g}_W} - 0.463\frac{\delta\Gamma_W}{\Gamma_W^{\text{SM}}} - 9.643\frac{\delta m_W}{m_W} - 1.487\bar{C}_{HW}. \quad (\text{A.32})$$

The SMEFT correction to the total Higgs width is finally estimated as

$$\frac{\delta\Gamma_h}{\Gamma_h^{\text{SM}}} = \sum_f \text{Br}_{h\rightarrow f}^{\text{SM,best}} \left[\frac{\delta\Gamma_{h\rightarrow f}}{\Gamma_{h\rightarrow f}^{\text{SM}}} \right], \quad (\text{A.33})$$

with $f = \{\bar{b}b, \bar{c}c, \tau^+\tau^-, \gamma\gamma, Z\gamma, gg, l^+\nu\bar{l}^-, \bar{u}d\bar{d}u, l^+\nu\bar{u}d + \text{h.c.}\}$ and the branching ratio values in table 6.

A.4 Top quark

To a very good approximation, the top quark decays exclusively to Wb . The SM width is

$$\Gamma_t^{\text{SM}} = \Gamma_{t\rightarrow W+b}^{\text{SM}} = \frac{G_F m_t^3}{8\sqrt{2}\pi} \left[(1-x_b^2)^2 + (1+x_b^2)x_W^2 - 2x_W^4 \right] \sqrt{\lambda(1, x_b^2, x_W^2)}, \quad (\text{A.34})$$

with $x_b = m_b/m_t$, $x_W = \hat{m}_W/m_t$ and $\lambda(a, b, c) = a^2 + b^2 + c^2 - 2ab - 2ac - 2bc$. The relative SMEFT correction is:

$$\begin{aligned} \frac{\delta\Gamma_t}{\Gamma_t^{\text{SM}}} &= \frac{\delta\Gamma_{t\rightarrow W+b}}{\Gamma_{t\rightarrow W+b}^{\text{SM}}} \\ &= 2\frac{\delta g_W}{\hat{g}_W} + 2(\bar{C}_{Hq}^{(3)})_{33} + \frac{6x_W}{(1-x_b^2)^2 + (1+x_b^2)x_W^2 - 2x_W^4} \left[x_W x_b (\text{Re}\bar{C}_{Hud})_{33} + \right. \\ &\quad \left. -\sqrt{2}(\text{Re}\bar{C}_{uW})_{33} (1-x_W^2 - x_b^2) + \sqrt{2}(\text{Re}\bar{C}_{dW})_{33} x_b (1+x_W^2 - x_b^2) \right] + \\ &\quad + 2 \left[\frac{(1-x_b^2)^2 - 3(1+x_b^2)x_W^2 + 2x_W^4}{\lambda(1, x_b^2, x_W^2)} - \frac{2(1-x_b^2)^2 + (1+x_b^2)^2 x_W^2}{(1-x_b^2)^2 + (1+x_b^2)x_W^2 - 2x_W^4} \right] \frac{\delta m_W}{\hat{m}_W}. \end{aligned} \quad (\text{A.35})$$

Decaying the W in final state does not lead to any additional contribution to $\delta\Gamma_t/\Gamma_t^{\text{SM}}$. This happens because the W is always on-shell, so its decay essentially factorizes out: using the narrow width approximation, one trivially has

$$\Gamma_t = \sum_f \Gamma_{t\rightarrow W+b} \cdot \text{Br}_{W\rightarrow f} = \Gamma_{t\rightarrow W+b}. \quad (\text{A.36})$$

Note that this conclusion only holds at the SM- \mathcal{L}_6 interference level, while at $\mathcal{O}(\Lambda^{-4})$ additional SMEFT corrections arise through contact vertices $(\bar{t}b)(\bar{\nu}l)$, $(\bar{t}b)(\bar{d}u)$.

B What's new in version 3.0

Here we briefly summarize the most significant updates and features introduced in SMEFTsim 3.0, compared to previous versions:

- The flavor assumptions `top` and `topU31` described in section 3.4 have been added.
- The flavor structure of all models has been generally improved. See section 3.5 for details.
- The treatment of propagator corrections described in section 6 has been implemented, enabling the estimate of linearized EFT corrections to the W, Z, H, t widths and to the W mass.
- The treatment of SM loop-generated Higgs interactions has been improved, particularly in the Higgs-gluon case. See section 5 for the general treatment and section 5.2 for a detailed comparison with previous implementations.
- All interaction vertices with up to 6 legs are now included in all UFO models. In the previous version, only 4-point functions were retained.
- The numerical values of the SM parameters have been updated, see table 16. The default value has been set to 0 for all Wilson coefficients.
- All complex Wilson coefficients are expressed in terms of their real and imaginary parts, rather than absolute values and phases.
- Individual interaction orders have been defined for each operator. Additionally, interaction orders `NPcpv`, `NPprop`, `NPfv` have been added, to provide more control on each class of EFT contributions. See section 8.2 for further details.
- In the UFO models, the SMEFT parameters have been organized in parameters blocks: `SMEFTcutoff` contains only Λ and `SMEFT` (`SMEFTcpv`) contain CP conserving (violating) Wilson coefficients. In the MFV models the flavor-violating ΔC_α parameters are contained in the additional `SMEFTFV` block. See section 8.1.
- The normalization of ΔG_F (`dGf`) has been modified. This is explicit in the `FeynRules` Lagrangian but does not have any consequence for the UFO models.

C Conversion tables between flavor assumptions

This appendix collects the results of section 3 and compares the flavor structure of the fermionic operators across the five setups considered. Tables 10–15 provide a dictionary between the different models: in order to translate between two flavor assumptions it is sufficient to exchange the corresponding expressions within each table block. All structures are given explicitly in terms of diagonal Yukawa matrices and of the CKM matrix V . In the `top` and `topU31` cases, Y_u and Y_d are 2×2 matrices and $V = \mathbb{1}$ is assumed.

general	$(C_{eH})_{pr}$		
U35, MFV, topU31	$C_{eH}(Y_l^{(d)})_{pr}$		
top	$(C_{eH})_{pp} \delta_{pr}$		
general	$(C_{uH})_{pr}$		
U35	$C_{uH}(Y_u^{(d)})_{pr}$		
MFV	$[C_{uH}^{(0)} Y_u^{(d)} + (\Delta^u C_{uH})(Y_u^{(d)})^3 + (\Delta^d C_{uH}) V(Y_d^{(d)})^2 V^\dagger Y_u^{(d)}]_{pr}$		
top, topU31	$C_{uH}(Y_u^{(d)})_{pr}$	$p, r = \{1, 2\},$	$C_{tH} \quad p = r = 3$
general	$(V^\dagger C_{dH})_{pr}$		
U35	$C_{dH}(Y_d^{(d)})_{pr}$		
MFV	$[C_{dH}^{(0)} Y_d^{(d)} + (\Delta^u C_{dH}) V^\dagger (Y_u^{(d)})^2 V Y_d^{(d)} + (\Delta^d C_{dH})(Y_d^{(d)})^3]_{pr}$		
top, topU31	$C_{dH}(Y_d^{(d)})_{pr}$	$p, r = \{1, 2\},$	$C_{bH} \quad p = r = 3$
general	$(C_{eW})_{pr}$		Same as $(C_{eH})_{pr}$.
general	$(C_{eB})_{pr}$		Same as $(C_{eH})_{pr}$.
general	$(C_{uW})_{pr}$		Same as $(C_{uH})_{pr}$.
general	$(C_{uB})_{pr}$		Same as $(C_{uH})_{pr}$.
general	$(C_{uG})_{pr}$		Same as $(C_{uH})_{pr}$.
general	$(V^\dagger C_{dW})_{pr}$		Same as $(V^\dagger C_{dH})_{pr}$.
general	$(V^\dagger C_{dB})_{pr}$		Same as $(V^\dagger C_{dH})_{pr}$.
general	$(V^\dagger C_{dG})_{pr}$		Same as $(V^\dagger C_{dH})_{pr}$.

Table 10. Conversion table among the 5 flavor assumptions considered for the operators in $\mathcal{L}_6^{(5)}, \mathcal{L}_6^{(6)}$.

general	$(C_{Hl}^{(1)})_{pr}$	
U35, MFV, topU3l	$C_{Hl}^{(1)} \delta_{pr}$	
top	$(C_{Hl}^{(1)})_{pp} \delta_{pr}$	
general	$(C_{Hl}^{(3)})_{pr}$	Same as $(C_{Hl}^{(1)})_{pr}$.
general	$(C_{He})_{pr}$	Same as $(C_{Hl}^{(1)})_{pr}$.
general	$(C_{Hq}^{(1)})_{pr}$	
U35	$C_{Hq}^{(1)} \delta_{pr}$	
MFV	$\left[C_{Hq}^{(1)(0)} \mathbb{1} + (\Delta^u C_{Hq}^{(1)}) (Y_u^{(d)})^2 + (\Delta^d C_{Hq}^{(1)}) V(Y_d^{(d)})^2 V^\dagger \right]_{pr}$	
top, topU3l	$C_{Hq}^{(1)} \delta_{pr} \quad p, r = \{1, 2\},$	$C_{Hq}^{(1)} \quad p = r = 3$
general	$(C_{Hq}^{(3)})_{pr}$	Same as $(C_{Hq}^{(1)})_{pr}$.
general	$(C_{Hu})_{pr}$	
U35	$C_{Hu} \delta_{pr}$	
MFV	$\left[C_{Hu}^{(0)} \mathbb{1} + (\Delta C_{Hu}) (Y_u^{(d)})^2 \right]_{pr}$	
top, topU3l	$C_{Hu} \delta_{pr} \quad p, r = \{1, 2\},$	$C_{Ht} \quad p = r = 3$
general	$(C_{Hd})_{pr}$	Same as $(C_{Hu})_{pr}$ with $Y_u^{(d)} \rightarrow Y_d^{(d)}, C_{Ht} \rightarrow C_{Hb}$.
general	$(C_{Hud})_{pr}$	
U35	$C_{Hud} \left[Y_u^{(d)} V Y_d^{(d)} \right]_{pr}$	
MFV	$C_{Hud}^{(0)} \left[Y_u^{(d)} V Y_d^{(d)} \right]_{pr}$	
top, topU3l	$C_{Hud} \left[Y_u^{(d)} Y_d^{(d)} \right]_{pr} \quad p, r = \{1, 2\},$	$C_{Htb} \quad p = r = 3$

Table 11. Conversion table among the 5 flavor assumptions considered, for the operators in $\mathcal{L}_6^{(7)}$.

general	$(C_{ll})_{prst}$	
U35, MFV, topU31	$C_{ll} \delta_{pr} \delta_{st} + C'_{ll} \delta_{pt} \delta_{sr}$	
top	$\frac{1}{2}(C_{ll})_{prst} (\delta_{pr} \delta_{st} + \delta_{pt} \delta_{sr}), \quad prst \in P_{ll}$	
general	$(C_{qq}^{(1)})_{prst}$	
U35	$C_{qq}^{(1)} \delta_{pr} \delta_{st} + C_{qq}^{(1)'} \delta_{pt} \delta_{sr}$	
MFV	$C_{qq}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{qq}^{(1)}) \left[(Y_u^{(d)})^2 \right]_{pr} \delta_{st} + (\Delta^d C_{qq}^{(1)}) \left[V(Y_d^{(d)})^2 V^\dagger \right]_{pr} \delta_{st}$ $+ C_{qq}^{(1) '(0)} \delta_{pt} \delta_{sr} + (\Delta^u C_{qq}^{(1)'}) \left[(Y_u^{(d)})^2 \right]_{pt} \delta_{sr} + (\Delta^d C_{qq}^{(1)'}) \left[V(Y_d^{(d)})^2 V^\dagger \right]_{pt} \delta_{sr}$	
top, topU31	$\left(C_{qq}^{(1,1)} - \frac{1}{6} C_{qq}^{(1,8)} \right) \delta_{pr} \delta_{st} + \frac{1}{4} \left(C_{qq}^{(1,8)} + 3C_{qq}^{(3,8)} \right) \delta_{pt} \delta_{sr} \quad p, r, s, t = \{1, 2\},$ $\left(\frac{1}{2} C_{Qq}^{(1,1)} - \frac{1}{12} C_{Qq}^{(1,8)} \right) \delta_{pr} \quad p, r = \{1, 2\}, s = t = 3,$ $\frac{1}{8} \left(C_{Qq}^{(1,8)} + 3C_{Qq}^{(3,8)} \right) \delta_{pt} \quad p, t = \{1, 2\}, s = r = 3,$ $C_{QQ}^{(1)} + \frac{1}{12} C_{QQ}^{(8)} \quad p = r = s = t = 3$	
general	$(C_{qq}^{(3)})_{prst}$	
U35	$C_{qq}^{(3)} \delta_{pr} \delta_{st} + C_{qq}^{(3)'} \delta_{pt} \delta_{sr}$	
MFV	$C_{qq}^{(3)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{qq}^{(3)}) \left[(Y_u^{(d)})^2 \right]_{pr} \delta_{st} + (\Delta^d C_{qq}^{(3)}) \left[V(Y_d^{(d)})^2 V^\dagger \right]_{pr} \delta_{st}$ $+ C_{qq}^{(3) '(0)} \delta_{pt} \delta_{sr} + (\Delta^u C_{qq}^{(3)'}) \left[(Y_u^{(d)})^2 \right]_{pt} \delta_{sr} + (\Delta^d C_{qq}^{(3)'}) \left[V(Y_d^{(d)})^2 V^\dagger \right]_{pt} \delta_{sr}$	
top, topU31	$\left(C_{qq}^{(3,1)} - \frac{1}{6} C_{qq}^{(3,8)} \right) \delta_{pr} \delta_{st} + \frac{1}{4} \left(C_{qq}^{(1,8)} - C_{qq}^{(3,8)} \right) \delta_{pt} \delta_{sr} \quad p, r, s, t = \{1, 2\},$ $\left(\frac{1}{2} C_{Qq}^{(3,1)} - \frac{1}{12} C_{Qq}^{(3,8)} \right) \delta_{pr} \quad p, r = \{1, 2\}, s = t = 3,$ $\frac{1}{8} \left(C_{Qq}^{(1,8)} - C_{Qq}^{(3,8)} \right) \delta_{pt} \quad p, t = \{1, 2\}, s = r = 3,$ $\frac{1}{4} C_{QQ}^{(8)} \quad p = r = s = t = 3$	
general	$(C_{lq}^{(1)})_{prst}$	
U35	$C_{lq}^{(1)} \delta_{pr} \delta_{st}$	
MFV	$C_{lq}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{lq}^{(1)}) \delta_{pr} \left[(Y_u^{(d)})^2 \right]_{st} + (\Delta^d C_{lq}^{(1)}) \delta_{pr} \left[V(Y_d^{(d)})^2 V^\dagger \right]_{st}$	
top	$(C_{lq}^{(1)})_{pp} \delta_{pr} \delta_{st} \quad s, t = \{1, 2\}, \quad (C_{lQ}^{(1)})_{pp} \delta_{pr} \quad s = t = 3$	
topU31	$C_{lq}^{(1)} \delta_{pr} \delta_{st} \quad s, t = \{1, 2\}, \quad C_{lQ}^{(1)} \delta_{pr} \quad s = t = 3$	
general	$(C_{lq}^{(3)})_{prst}$	Same as $(C_{lq}^{(1)})_{prst}$.

Table 12. Conversion table among the 5 flavor assumptions considered, for the operators in $\mathcal{L}_6^{(8a)}$. The set P_{ll} is defined in eq. (3.70).

general	$(C_{ee})_{prst}$	
U35, MFV, topU31	$\frac{1}{2}C_{ee}(\delta_{pr}\delta_{st} + \delta_{pt}\delta_{sr})$	
top	$\frac{1}{4}(C_{ee})_{prst}(\delta_{pr}\delta_{st} + 3\delta_{pt}\delta_{sr}),$	$prst \in P_{ee}$
general	$(C_{uu})_{prst}$	
U35	$C_{uu}\delta_{pr}\delta_{st} + C'_{uu}\delta_{pt}\delta_{sr}$	
MFV	$C_{uu}^{(0)}\delta_{pr}\delta_{st} + (\Delta C_{uu})\left[(Y_u^{(d)})^2\right]_{pr}\delta_{st} + C_{uu}^{(0)'}\delta_{pt}\delta_{sr} + (\Delta C'_{uu})\left[(Y_u^{(d)})^2\right]_{pt}\delta_{sr}$	
top	$\left(C_{uu}^{(1)} - \frac{1}{6}C_{uu}^{(8)}\right)\delta_{pr}\delta_{st} + \frac{1}{2}C_{uu}^{(8)}\delta_{pt}\delta_{sr}$	$p, r, s, t = \{1, 2\},$
	$\left(\frac{1}{2}C_{tu}^{(1)} - \frac{1}{12}C_{tu}^{(8)}\right)\delta_{pr}$	$p, r = \{1, 2\}, s = t = 3,$
	$\frac{1}{4}C_{tu}^{(8)}\delta_{pt}$	$p, t = \{1, 2\}, s = r = 3,$
	C_{tt}	$p = r = s = t = 3$
general	$(C_{dd})_{prst}$	Same as $(C_{uu})_{prst}$ with $Y_u^{(d)} \rightarrow Y_d^{(d)},$ $C_{tu}^{(1),(8)} \rightarrow C_{bd}^{(1),(8)}, C_{tt} \rightarrow C_{bb}.$
general	$(C_{eu})_{prst}$	
U35	$C_{eu}\delta_{pr}\delta_{st}$	
MFV	$C_{eu}^{(0)}\delta_{pr}\delta_{st} + (\Delta C_{eu})\delta_{pr}\left[(Y_u^{(d)})^2\right]_{st}$	
top	$(C_{eu})_{pp}\delta_{pr}\delta_{st}$	$s, t = \{1, 2\},$ $(C_{et})_{pp}\delta_{pr}$ $s = t = 3,$
topU31	$C_{eu}\delta_{pr}\delta_{st}$	$s, t = \{1, 2\},$ $C_{et}\delta_{pr}$ $s = t = 3$
general	$(C_{ed})_{prst}$	Same as $(C_{eu})_{prst}$ with $Y_u^{(d)} \rightarrow Y_d^{(d)}, C_{et} \rightarrow C_{eb}.$
general	$(C_{ud}^{(1)})_{prst}$	
U35	$C_{ud}^{(1)}\delta_{pr}\delta_{st}$	
MFV	$C_{ud}^{(1)(0)}\delta_{pr}\delta_{st} + (\Delta^u C_{ud}^{(1)})\left[(Y_u^{(d)})^2\right]_{pr}\delta_{st} + (\Delta^d C_{ud}^{(1)})\delta_{pr}\left[(Y_d^{(d)})^2\right]_{st}$	
top, topU31	$C_{ud}^{(1)}\delta_{pr}\delta_{st}$	$p, r, s, t = \{1, 2\},$ $C_{tb}^{(1)}$ $p = r = s = t = 3,$
	$C_{ub}^{(1)}\delta_{pr}$	$p, r = \{1, 2\}, s = t = 3,$ $C_{td}^{(1)}\delta_{st}$ $s, t = \{1, 2\}, p = r = 3$
	$C_{utbd}^{(1)}\left[Y_u^{(d)}Y_d^{(d)}\right]_{pt}$	$p, t = \{1, 2\}, s = r = 3,$
	$C_{utbd}^{(1)\star}\left[Y_d^{(d)}Y_u^{(d)}\right]_{sr}$	$s, r = \{1, 2\}, p = t = 3,$
general	$(C_{ud}^{(8)})_{prst}$	Same as $(C_{ud}^{(1)})_{prst}.$

Table 13. Conversion table among the 5 flavor assumptions considered, for the operators in $\mathcal{L}_6^{(8b)}$. The set P_{ee} is defined in eq. (3.91).

general	$(C_{le})_{prst}$		
U35, MFV, topU31	$C_{le} \delta_{pr} \delta_{st}$		
top	$(C_{le})_{prst}$	$prst \in P_{le}$	
general	$(C_{lu})_{prst}$		
U35	$C_{lu} \delta_{pr} \delta_{st}$		
MFV	$C_{lu}^{(0)} \delta_{pr} \delta_{st} + (\Delta C_{lu}) \delta_{pr} [(Y_u^{(d)})^2]_{st}$		
top	$(C_{lu})_{pp} \delta_{pr} \delta_{st}$	$s, t = \{1, 2\},$	$(C_{lt})_{pp} \delta_{pr} \quad s = t = 3,$
topU31	$C_{lu} \delta_{pr} \delta_{st}$	$s, t = \{1, 2\},$	$C_{lt} \delta_{pr} \quad s = t = 3$
general	$(C_{ld})_{prst}$		Same as $(C_{lu})_{prst}$ with $Y_u^{(d)} \rightarrow Y_d^{(d)}, C_{lt} \rightarrow C_{lb}.$
general	$(C_{qe})_{prst}$		
U35	$C_{qe} \delta_{pr} \delta_{st}$		
MFV	$C_{qe}^{(0)} \delta_{pr} \delta_{st} + (\Delta^u C_{qe}) [(Y_u^{(d)})^2]_{pr} \delta_{st} + (\Delta^d C_{qe}) [V(Y_d^{(d)})^2 V^\dagger]_{pr} \delta_{st}$		
top	$(C_{qe})_{pp} \delta_{pr} \delta_{st}$	$p, r = \{1, 2\},$	$(C_{Qe})_{pp} \delta_{st} \quad p = r = 3,$
topU31	$C_{qe} \delta_{pr} \delta_{st}$	$p, r = \{1, 2\},$	$C_{Qe} \delta_{st} \quad p = r = 3$
general	$(C_{qu}^{(1)})_{prst}$		
U35	$C_{qu}^{(1)} \delta_{pr} \delta_{st}$		
MFV	$C_{qu}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta_1^u C_{qu}^{(1)}) [(Y_u^{(d)})^2]_{pr} \delta_{st} + (\Delta_1^d C_{qu}^{(1)}) [V(Y_d^{(d)})^2 V^\dagger]_{pr} \delta_{st}$ $+ (\Delta_2 C_{qu}^{(1)}) \delta_{pr} [(Y_u^{(d)})^2]_{st} + C_{qu}^{(1) \prime(0)} (Y_u^{(d)})_{pt} (Y_u^{(d)})_{sr}$		
top, topU31	$C_{qu}^{(1)} \delta_{pr} \delta_{st}$	$p, r, s, t = \{1, 2\},$	$C_{Qt}^{(1)} \quad p = r = s = t = 3,$
	$C_{qt}^{(1)} \delta_{pr}$	$p, r = \{1, 2\}, s = t = 3,$	$C_{Qu}^{(1)} \delta_{st} \quad s, t = \{1, 2\}, p = r = 3$
	$C_{qQtu}^{(1)} [Y_u^{(d)}]_{pt}$	$p, t = \{1, 2\}, s = r = 3$	$C_{qQtu}^{(1)*} [Y_u^{(d)}]_{sr} \quad s, r = \{1, 2\}, p = t = 3$
general	$(C_{qu}^{(8)})_{prst}$		Same as $(C_{qu}^{(1)})_{prst}.$
general	$(C_{qd}^{(1)})_{prst}$		
U35	$C_{qd}^{(1)} \delta_{pr} \delta_{st}$		
MFV	$C_{qd}^{(1)(0)} \delta_{pr} \delta_{st} + (\Delta_1^u C_{qd}^{(1)}) [(Y_u^{(d)})^2]_{pr} \delta_{st} + (\Delta_1^d C_{qd}^{(1)}) [V(Y_d^{(d)})^2 V^\dagger]_{pr} \delta_{st}$ $+ (\Delta_2 C_{qd}^{(1)}) \delta_{pr} [(Y_d^{(d)})^2]_{st} + C_{qd}^{(1) \prime(0)} [V Y_d^{(d)}]_{pt} [Y_d^{(d)} V^\dagger]_{sr}$		
top, topU31	$C_{qd}^{(1)} \delta_{pr} \delta_{st}$	$p, r, s, t = \{1, 2\},$	$C_{Qb}^{(1)} \quad p = r = s = t = 3,$
	$C_{qb}^{(1)} \delta_{pr}$	$p, r = \{1, 2\}, s = t = 3,$	$C_{Qd}^{(1)} \delta_{st} \quad s, t = \{1, 2\}, p = r = 3$
general	$(C_{qd}^{(8)})_{prst}$		Same as $(C_{qd}^{(1)})_{prst}.$

Table 14. Conversion table among the 5 flavor assumptions considered, for the operators in $\mathcal{L}_6^{(8c)}$. The set P_{le} is defined in eq. (3.93).

general	$(C_{ledq})_{prst}$			
U35	$C_{ledq} [Y_l^{(d)}]_{pr} [Y_d^{(d)} V^\dagger]_{st}$			
MFV	$C_{ledq}^{(0)} [Y_l^{(d)}]_{pr} [Y_d^{(d)} V^\dagger]_{st} + (\Delta^u C_{ledq}) [Y_l^{(d)}]_{pr} [Y_d^{(d)} V^\dagger (Y_u^{(d)})^2]_{st}$ $+ (\Delta^d C_{ledq}) [Y_l^{(d)}]_{pr} [(Y_d^{(d)})^3 V^\dagger]_{st}$			
top	$(C_{ledq})_{pp} \delta_{pr} [Y_d^{(d)}]_{st}$	$s, t = \{1, 2\},$	$(C_{lebQ})_{pp} \delta_{pr}$	$s = t = 3$
topU31	$C_{ledq} [Y_l^{(d)}]_{pr} [Y_d^{(d)}]_{st}$	$s, t = \{1, 2\},$	$C_{lebQ} [Y_l^{(d)}]_{pr}$	$s = t = 3$
general	$(C_{lequ}^{(1)})_{prst}$			
U35	$C_{lequ}^{(1)} [Y_l^{(d)}]_{pr} [Y_u^{(d)}]_{st}$			
MFV	$C_{lequ}^{(1)(0)} [Y_l^{(d)}]_{pr} [Y_u^{(d)}]_{st} + (\Delta^u C_{lequ}^{(1)}) [Y_l^{(d)}]_{pr} [(Y_u^{(d)})^3]_{st}$ $+ (\Delta^d C_{lequ}^{(1)}) [Y_l^{(d)}]_{pr} [V (Y_d^{(d)})^2 V^\dagger Y_u^{(d)}]_{st}$			
top	$(C_{lequ}^{(1)})_{pp} \delta_{pr} [Y_u^{(d)}]_{st}$	$s, t = \{1, 2\},$	$(C_{leQt}^{(1)})_{pp} \delta_{pr}$	$s = t = 3$
topU31	$C_{lequ}^{(1)} [Y_l^{(d)}]_{pr} [Y_u^{(d)}]_{st}$	$s, t = \{1, 2\},$	$C_{leQt}^{(1)} [Y_l^{(d)}]_{pr}$	$s = t = 3$
general	$(C_{lequ}^{(3)})_{prst}$		Same as $(C_{lequ}^{(1)})_{prst}$.	
general	$(C_{quqd}^{(1)})_{prst}$			
U35, MFV	$C_{quqd}^{(1)} [Y_u^{(d)}]_{pr} [V Y_d^{(d)}]_{st} + C_{quqd}^{(1)'} [Y_u^{(d)}]_{sr} [V Y_d^{(d)}]_{pt}$			
top, topU31	$C_{quqd}^{(1)} [Y_u^{(d)}]_{pr} [Y_d^{(d)}]_{st} + C_{quqd}^{(1)'} [Y_u^{(d)}]_{sr} [Y_d^{(d)}]_{pt}$	$p, r, s, t = \{1, 2\},$		
	$C_{QtQb}^{(1)}$		$p = r = s = t = 3,$	
	$C_{quQb}^{(1)} [Y_u^{(d)}]_{pr}$		$p, r = \{1, 2\}, s = t = 3,$	
	$C_{Qtqd}^{(1)} [Y_d^{(d)}]_{st}$		$s, t = \{1, 2\}, p = r = 3,$	
	$C_{Quqb}^{(1)} [Y_u^{(d)}]_{sr}$		$s, r = \{1, 2\}, p = t = 3,$	
	$C_{qtQd}^{(1)} [Y_d^{(d)}]_{pt}$		$p, t = \{1, 2\}, s = r = 3$	
general	$(C_{quqd}^{(8)})_{prst}$		Same as $(C_{quqd}^{(1)})_{prst}$.	

Table 15. Conversion table among the 5 flavor assumptions considered, for the operators in $\mathcal{L}_6^{(8d)}$.

D Parameter definitions in the code implementation

This appendix provides tables to facilitate the interpretation of the `FeynRules` and `UFO` implementations in terms of the theory discussion in the main text.

Table 16 lists the external parameters that are defined in all `SMEFTsim` models, specifying the corresponding code name and default numerical value. Table 17 shows the nomenclature used for the Wilson coefficients of the bosonic operators and for the shift quantities defined in section 4.1. Tables 18–23 do the same for the Wilson coefficients of fermionic operators, for each flavor assumption. As a common rationale, primes are replaced by 1 in the code name and real and imaginary parts are specified by with `Re`, `Im` suffixes. If needed, flavor indices are fully specified and appended at the very end of the code names. In the `MFV` models, the coefficients $\Delta_n^q(C_\alpha)$ are denoted `Delta[n][q]c[a]`.

Although the correspondence between parameters names is most often direct, some notational changes were necessary, particularly in the `top` and `topU31` implementations. Most notably the lowercase q has been replaced with `j` in all the parameters' and operators' names, as the q/Q distinction between light and heavy quark fields is problematic for non-case-sensitive interfaces. Analogously, the coefficient C_{Hb} is denoted as `cHbq` to avoid conflict with `cHB`, while C_{bB} is denoted as `cbBB`, distinct from `cbb`. The internal parameter C_{tH} is denoted as `ctHH` to avoid conflict with the cosine of the weak angle `cth`.

Common parameters defined in SMEFTsim							
parameter	UFO	default value		parameter	UFO	default value	
G_F	Gf	1.1663787×10^{-5} GeV ⁻²	[82, 118]	$\alpha_s(m_Z)$	aS	0.1179	[82]
$\alpha_{em}(m_Z)$ (*)	aEW	1/127.95	[82, 118]	m_W (**)	MW	80.387 GeV	[119]
m_d	MD	4.67×10^{-3} GeV	[82]	$y_d \hat{v}/\sqrt{2}$	ymdo	4.67×10^{-3} GeV	[82]
m_s	MS	0.093 GeV	[82]	$y_s \hat{v}/\sqrt{2}$	yms	0.093 GeV	[82]
m_b	MB	4.18 GeV	[82]	$y_b \hat{v}/\sqrt{2}$	ymb	4.18 GeV	[82]
m_u	MU	2.16×10^{-3} GeV	[82]	$y_u \hat{v}/\sqrt{2}$	ymup	2.16×10^{-3} GeV	[82]
m_c	MC	1.27 GeV	[82]	$y_c \hat{v}/\sqrt{2}$	ymc	1.27 GeV	[82]
m_t	MT	172.76 GeV	[82]	$y_t \hat{v}/\sqrt{2}$	ymt	172.76 GeV	[82]
m_e	Me	5.11×10^{-4} GeV	[82]	$y_e \hat{v}/\sqrt{2}$	yme	5.11×10^{-4} GeV	[82]
m_μ	MMU	0.10566 GeV	[82]	$y_\mu \hat{v}/\sqrt{2}$	ymm	0.10566 GeV	[82]
m_τ	MTA	1.777 GeV	[82]	$y_\tau \hat{v}/\sqrt{2}$	ymtau	1.777 GeV	[82]
m_Z	MZ	91.1876 GeV	[82, 118, 120]	m_h	MH	125.09 GeV	[82]
$\Gamma_Z^{SM,best}$	WZ	2.4952 GeV	[82]	$\Gamma_h^{SM,best}$	WH	4.07×10^{-3} GeV	[113]
$\Gamma_W^{SM,best}$	WW	2.085 GeV	[82]	$\Gamma_t^{SM,best}$	WT	1.33 GeV	[105]
λ_{CKM}	CKMlambda	0.22650	[82]	A	CKMA	0.790	[82]
ρ	CKMrho	0.141	[82]	η	CKMeta	0.357	[82]
Λ	LambdaSMEFT	1 TeV					

(*) only in models with $\{\alpha_{em}, m_Z, G_F\}$ inputs. (**) only in models with $\{m_W, m_Z, G_F\}$ inputs.

Table 16. Common external parameters defined in all models, with the corresponding name in the code and default numerical value. The latter have been updated compared to previous SMEFTsim versions, according to the references indicated.

Bosonic SMEFT parameters in SMEFTsim								
$\mathcal{L}_6^{(1)}$	C_G	cG	$C_{\tilde{G}}$	cGtil	C_W	cW	$C_{\tilde{W}}$	cWtil
$\mathcal{L}_6^{(2,3)}$	C_H	cH	$C_{H\Box}$	cHbox	C_{HD}	cHDD		
$\mathcal{L}_6^{(4)}$	C_{HG}	cHG	C_{HW}	cHW	C_{HB}	cHB	C_{HWB}	cHWB
	$C_{H\tilde{G}}$	cHGtil	$C_{H\tilde{W}}$	cHWtil	$C_{H\tilde{B}}$	cHBtil	$C_{H\tilde{W}B}$	cHWBtil
Shift parameters in SMEFTsim								
Δm_Z^2	dMZ2		Δm_h^2	dMH2	ΔG_F	dGf	$\Delta \kappa_H$	dkH
$\delta \Gamma_Z$	dWZ		$\delta \Gamma_W$	dWW	$\delta \Gamma_h$	dWH	$\delta \Gamma_t$	dWT
$\delta g_1/\hat{g}_1$	dg1		$\delta g_W/\hat{g}_W$	dgw	δm_W	dMW		

Table 17. Upper block: Wilson coefficients for the 15 bosonic operators. They are common to all flavor versions except MFV, where the CP violating $C_{\tilde{G}}, C_{\tilde{W}}, C_{H\tilde{G}}, C_{H\tilde{W}}, C_{H\tilde{B}}, C_{H\tilde{W}B}$ are not defined. Lower block: shift parameters defined in SMEFTsim, see section 4.1 for details. δm_W is defined only in models with the $\{\alpha_{em}, m_Z, G_F\}$ input scheme.

Fermionic SMEFT parameters in SMEFTsim general						
class	parameter	UFO	[pr(st)]	parameter	UFO	[pr(st)]
$\mathcal{L}_6^{(5)}$	$\text{Re}(C_{eH})_{pr}$	ceHRe [pr]	[2f-NH-R]	$\text{Im}(C_{eH})_{pr}$	ceHIm [pr]	[2f-NH-I]
	$\text{Re}(C_{uH})_{pr}$	cdHRe [pr]	[2f-NH-R]	$\text{Im}(C_{uH})_{pr}$	cdHIm [pr]	[2f-NH-I]
	$\text{Re}(C_{dH})_{pr}$	cdHRe [pr]	[2f-NH-R]	$\text{Im}(C_{dH})_{pr}$	cdHIm [pr]	[2f-NH-I]
$\mathcal{L}_6^{(6)}$	$\text{Re}(C_{eW})_{pr}$	ceWRe [pr]	[2f-NH-R]	$\text{Im}(C_{eW})_{pr}$	ceWIm [pr]	[2f-NH-I]
	$\text{Re}(C_{eB})_{pr}$	ceBRe [pr]	[2f-NH-R]	$\text{Im}(C_{eB})_{pr}$	ceBIm [pr]	[2f-NH-I]
	$\text{Re}(C_{uG})_{pr}$	cuGRe [pr]	[2f-NH-R]	$\text{Im}(C_{uG})_{pr}$	cuGIm [pr]	[2f-NH-I]
	$\text{Re}(C_{uW})_{pr}$	cuWRe [pr]	[2f-NH-R]	$\text{Im}(C_{uW})_{pr}$	cuWIm [pr]	[2f-NH-I]
	$\text{Re}(C_{uB})_{pr}$	cuBRe [pr]	[2f-NH-R]	$\text{Im}(C_{uB})_{pr}$	cuBIm [pr]	[2f-NH-I]
	$\text{Re}(C_{dG})_{pr}$	cdGRe [pr]	[2f-NH-R]	$\text{Im}(C_{dG})_{pr}$	cdGIm [pr]	[2f-NH-I]
	$\text{Re}(C_{dW})_{pr}$	cdWRe [pr]	[2f-NH-R]	$\text{Im}(C_{dW})_{pr}$	cdWIm [pr]	[2f-NH-I]
	$\text{Re}(C_{dB})_{pr}$	cdBRe [pr]	[2f-NH-R]	$\text{Im}(C_{dB})_{pr}$	cdBIm [pr]	[2f-NH-I]
$\mathcal{L}_6^{(7)}$	$\text{Re}(C_{H1}^{(1)})_{pr}$	ch11Re [pr]	[2f-H-R]	$\text{Im}(C_{H1}^{(1)})_{pr}$	ch11Im [pr]	[2f-H-I]
	$\text{Re}(C_{H1}^{(3)})_{pr}$	ch13Re [pr]	[2f-H-R]	$\text{Im}(C_{H1}^{(3)})_{pr}$	ch13Im [pr]	[2f-H-I]
	$\text{Re}(C_{He})_{pr}$	cHeRe [pr]	[2f-H-R]	$\text{Im}(C_{He})_{pr}$	cHeIm [pr]	[2f-H-I]
	$\text{Re}(C_{Hq}^{(1)})_{pr}$	chq1Re [pr]	[2f-H-R]	$\text{Im}(C_{Hq}^{(1)})_{pr}$	chq1Im [pr]	[2f-H-I]
	$\text{Re}(C_{Hq}^{(3)})_{pr}$	chq3Re [pr]	[2f-H-R]	$\text{Im}(C_{Hq}^{(3)})_{pr}$	chq3Im [pr]	[2f-H-I]
	$\text{Re}(C_{Hu})_{pr}$	chURe [pr]	[2f-H-R]	$\text{Im}(C_{Hu})_{pr}$	chUIm [pr]	[2f-H-I]
	$\text{Re}(C_{Hd})_{pr}$	chDRe [pr]	[2f-H-R]	$\text{Im}(C_{Hd})_{pr}$	chDIm [pr]	[2f-H-I]
	$\text{Re}(C_{Hud})_{pr}$	chudRe [pr]	[2f-NH-R]	$\text{Im}(C_{Hud})_{pr}$	chudIm [pr]	[2f-NH-I]
$\mathcal{L}_6^{(8a)}$	$\text{Re}(C_{ll})_{prst}$	cllRe [prst]	[4f-H-S-R]	$\text{Im}(C_{ll})_{prst}$	cllIm [prst]	[4f-H-S-I]
	$\text{Re}(C_{qq}^{(1)})_{prst}$	cqq1Re [prst]	[4f-H-S-R]	$\text{Im}(C_{qq}^{(1)})_{prst}$	cqq1Im [prst]	[4f-H-S-I]
	$\text{Re}(C_{qq}^{(3)})_{prst}$	cqq3Re [prst]	[4f-H-S-R]	$\text{Im}(C_{qq}^{(3)})_{prst}$	cqq3Im [prst]	[4f-H-S-I]
	$\text{Re}(C_{lq}^{(1)})_{prst}$	clq1Re [prst]	[4f-H-R]	$\text{Im}(C_{lq}^{(1)})_{prst}$	clq1Im [prst]	[4f-H-I]
	$\text{Re}(C_{lq}^{(3)})_{prst}$	clq3Re [prst]	[4f-H-R]	$\text{Im}(C_{lq}^{(3)})_{prst}$	clq3Im [prst]	[4f-H-I]
$\mathcal{L}_6^{(8b)}$	$\text{Re}(C_{ee})_{prst}$	ceeRe [prst]	[4f-ee-R]	$\text{Im}(C_{ee})_{prst}$	ceeIm [prst]	[4f-ee-I]
	$\text{Re}(C_{uu})_{prst}$	cuuRe [prst]	[4f-H-S-R]	$\text{Im}(C_{uu})_{prst}$	cuuIm [prst]	[4f-H-S-I]
	$\text{Re}(C_{dd})_{prst}$	cddRe [prst]	[4f-H-S-R]	$\text{Im}(C_{dd})_{prst}$	cddIm [prst]	[4f-H-S-I]
	$\text{Re}(C_{eu})_{prst}$	ceuRe [prst]	[4f-H-R]	$\text{Im}(C_{eu})_{prst}$	ceuIm [prst]	[4f-H-I]
	$\text{Re}(C_{ed})_{prst}$	cedRe [prst]	[4f-H-R]	$\text{Im}(C_{ed})_{prst}$	cedIm [prst]	[4f-H-I]
	$\text{Re}(C_{ud}^{(1)})_{prst}$	cud1Re [prst]	[4f-H-R]	$\text{Im}(C_{ud}^{(1)})_{prst}$	cud1Im [prst]	[4f-H-I]
	$\text{Re}(C_{ud}^{(8)})_{prst}$	cud8Re [prst]	[4f-H-R]	$\text{Im}(C_{ud}^{(8)})_{prst}$	cud8Im [prst]	[4f-H-I]
$\mathcal{L}_6^{(8c)}$	$\text{Re}(C_{le})_{prst}$	cleRe [prst]	[4f-H-R]	$\text{Im}(C_{le})_{prst}$	cleIm [prst]	[4f-H-I]
	$\text{Re}(C_{lu})_{prst}$	cluRe [prst]	[4f-H-R]	$\text{Im}(C_{lu})_{prst}$	cluIm [prst]	[4f-H-I]
	$\text{Re}(C_{ld})_{prst}$	clDRe [prst]	[4f-H-R]	$\text{Im}(C_{ld})_{prst}$	clDIm [prst]	[4f-H-I]
	$\text{Re}(C_{qe})_{prst}$	cqeRe [prst]	[4f-H-R]	$\text{Im}(C_{qe})_{prst}$	cqeIm [prst]	[4f-H-I]
	$\text{Re}(C_{qu}^{(1)})_{prst}$	cqu1Re [prst]	[4f-H-R]	$\text{Im}(C_{qu}^{(1)})_{prst}$	cqu1Im [prst]	[4f-H-I]
	$\text{Re}(C_{qu}^{(8)})_{prst}$	cqu8Re [prst]	[4f-H-R]	$\text{Im}(C_{qu}^{(8)})_{prst}$	cqu8Im [prst]	[4f-H-I]
	$\text{Re}(C_{qd}^{(1)})_{prst}$	cqd1Re [prst]	[4f-H-R]	$\text{Im}(C_{qd}^{(1)})_{prst}$	cqd1Im [prst]	[4f-H-I]
	$\text{Re}(C_{qd}^{(8)})_{prst}$	cqd8Re [prst]	[4f-H-R]	$\text{Im}(C_{qd}^{(8)})_{prst}$	cqd8Im [prst]	[4f-H-I]
$\mathcal{L}_6^{(8d)}$	$\text{Re}(C_{ledq})_{prst}$	clEdqRe [prst]	[4f-NH-R]	$\text{Im}(C_{ledq})_{prst}$	clEdqIm [prst]	[4f-NH-I]
	$\text{Re}(C_{quqd}^{(1)})_{prst}$	cquqd1Re [prst]	[4f-NH-R]	$\text{Im}(C_{quqd}^{(1)})_{prst}$	cquqd1Im [prst]	[4f-NH-I]
	$\text{Re}(C_{quqd}^{(8)})_{prst}$	cquqd8Re [prst]	[4f-NH-R]	$\text{Im}(C_{quqd}^{(8)})_{prst}$	cquqd8Im [prst]	[4f-NH-I]
	$\text{Re}(C_{lequ}^{(1)})_{prst}$	clequ1Re [prst]	[4f-NH-R]	$\text{Im}(C_{lequ}^{(1)})_{prst}$	clequ1Im [prst]	[4f-NH-I]
	$\text{Re}(C_{lequ}^{(3)})_{prst}$	clequ3Re [prst]	[4f-NH-R]	$\text{Im}(C_{lequ}^{(3)})_{prst}$	clequ3Im [prst]	[4f-NH-I]

Table 18. The 2484 independent parameters in $\mathcal{L}_6^{(5,6,7,8)}$ defined in the `general` flavor model, see section 3.1. The indices strings [pr], [prst] take values in the sets indicated, that are defined in table 19.

[2f-NH-R]	11, 22, 33, 12, 13, 23, 21, 31, 32
[2f-NH-I]	11, 22, 33, 12, 13, 23, 21, 31, 32
[2f-H-R]	11, 22, 33, 12, 13, 23
[2f-H-I]	12, 13, 23
[4f-ee-R]	1111, 1122, 1133, 2222, 2233, 3333, 1112, 1113, 1123, 1212, 1213, 1222, 1232, 1233, 1313, 1322, 1323, 1333, 2223, 2323, 3323
[4f-ee-I]	1112, 1113, 1123, 1212, 1213, 1222, 1232, 1233, 1313, 1322, 1323, 1333, 2223, 2323, 3323
[4f-H-S-R]	1111, 1122, 1133, 2222, 2233, 3333, 1221, 1331, 2332, 1112, 1113, 1123, 1212, 1213, 1222, 1232, 1233, 1313, 1322, 1323, 1333, 2223, 2323, 3323, 1231, 1223, 1332
[4f-H-S-I]	1112, 1113, 1123, 1212, 1213, 1222, 1232, 1233, 1313, 1322, 1323, 1333, 2223, 2323, 3323, 1231, 1223, 1332
[4f-H-R]	1111, 1122, 1133, 2222, 2233, 3333, 1221, 1331, 2332, 2211, 3311, 3322, 1112, 1113, 1123, 1212, 1213, 1222, 1232, 1233, 1313, 1322, 1323, 1333, 2223, 2323, 3323, 1231, 1223, 1332, 1211, 1311, 1312, 1321, 2212, 2213, 2311, 2312, 2313, 2321, 2322, 2331, 2333, 3312, 3313
[4f-H-I]	1112, 1113, 1123, 1212, 1213, 1222, 1232, 1233, 1313, 1322, 1323, 1333, 2223, 2323, 3323, 1231, 1223, 1332, 1211, 1311, 1312, 1321, 2212, 2213, 2311, 2312, 2313, 2321, 2322, 2331, 2333, 3312, 3313, 2211, 3311, 3322

Table 19. Sets of indices implemented for each category of fermionic operator in the `general` model.

Fermionic SMEFT parameters in SMEFTsim U35							
$\mathcal{L}_6^{(5)}$	Re C_{eH}	ceHRe	Re C_{uH}	cuHRe	Re C_{dH}	cdHRe	
	Im C_{eH}	ceHIm	Im C_{uH}	cuHIm	Im C_{dH}	cdHIm	
$\mathcal{L}_6^{(6)}$	Re C_{eW}	ceWRe	Re C_{eB}	ceBRe	Re C_{uG}	cuGRe	Re C_{uW} cuWRe
	Im C_{eW}	ceWIm	Im C_{eB}	ceBIm	Im C_{uG}	cuGIm	Im C_{uW} cuWIm
	Re C_{uB}	cuBRe	Re C_{dG}	cdGRe	Re C_{dW}	cdWRe	Re C_{dB} cdBRe
	Im C_{uB}	cuBIm	Im C_{dG}	cdGIm	Im C_{dW}	cdWIm	Im C_{dB} cdBIm
$\mathcal{L}_6^{(7)}$	$C_{Hl}^{(1)}$	cHl1	$C_{Hq}^{(1)}$	cHq1	C_{Hu}	cHu	C_{He} cHe
	$C_{Hl}^{(3)}$	cHl3	$C_{Hq}^{(3)}$	cHq3	C_{Hd}	cHd	
	Re C_{Hud}	cHudRe	Im C_{Hud}	cHudIm			
$\mathcal{L}_6^{(8a)}$	C_{ll}	c11	$C_{qq}^{(1)}$	cqq1	$C_{qq}^{(3)}$	cqq3	$C_{lq}^{(1)}$ clq1
	C'_{ll}	c111	$C_{qq}^{(1)'} $	cqq11	$C_{qq}^{(3)'} $	cqq31	$C_{lq}^{(3)}$ clq3
$\mathcal{L}_6^{(8b)}$	C_{uu}	cuu	C_{dd}	cdd	C_{eu}	ceu	$C_{ud}^{(1)}$ cud1
	C'_{uu}	cuu1	C'_{dd}	cdd1	C_{ed}	ced	$C_{ud}^{(8)}$ cud8
	C_{ee}	cee					
$\mathcal{L}_6^{(8c)}$	C_{le}	c1e	C_{lu}	c1u	$C_{qu}^{(1)}$	cqu1	$C_{qd}^{(1)}$ cqd1
	C_{qe}	cqe	C_{ld}	c1d	$C_{qu}^{(8)}$	cqu8	$C_{qd}^{(8)}$ cqd8
$\mathcal{L}_6^{(8d)}$	Re $C_{quqd}^{(1)}$	cquqd1Re	Re $C_{quqd}^{(8)}$	cquqd8Re	Im $C_{quqd}^{(1)}$	cquqd1Im	Im $C_{quqd}^{(8)}$ cquqd8Im
	Re $C_{quqd}^{(1)'}$	cquqd11Re	Re $C_{quqd}^{(8)'}$	cquqd81Re	Im $C_{quqd}^{(1)'}$	cquqd11Im	Im $C_{quqd}^{(8)'}$ cquqd81Im
	Re $C_{lequ}^{(1)}$	c1equ1Re	Re $C_{lequ}^{(3)}$	c1equ3Re	Im $C_{lequ}^{(1)}$	c1equ1Im	Im $C_{lequ}^{(3)}$ c1equ3Im
	Re C_{ledq}	c1edqRe	Im C_{ledq}	c1edqIm			

Table 20. The 70 independent parameters in $\mathcal{L}_6^{(5,6,7,8)}$ defined in the U35 model, see section 3.2.

Fermionic SMEFT parameters in SMEFTsim MFV										
$\mathcal{L}_6^{(5)}$	C_{eH} $(\Delta^u C_{uH})$	ceH DeltaucH	$C_{uH}^{(0)}$ $(\Delta^d C_{uH})$	cuH0 DeltadcuH	$C_{dH}^{(0)}$ $(\Delta^u C_{dH})$	cdH0 DeltaucdH	$(\Delta^d C_{dH})$	DeltadcdH		
$\mathcal{L}_6^{(6)}$	C_{eW} $C_{uB}^{(0)}$ $(\Delta^u C_{uG})$ $(\Delta^u C_{uW})$ $(\Delta^u C_{uB})$	ceW cuB0 DeltaucG DeltaucW DeltaucB	C_{eB} $C_{dG}^{(0)}$ $(\Delta^d C_{uG})$ $(\Delta^d C_{uW})$ $(\Delta^d C_{uB})$	ceB cdG0 DeltadcuG DeltadcuW DeltadcuB	$C_{uG}^{(0)}$ $C_{dW}^{(0)}$ $(\Delta^u C_{dG})$ $(\Delta^u C_{dW})$ $(\Delta^u C_{dB})$	cuG0 cdW0 DeltaucdG DeltaucdW DeltaucdB	$C_{uW}^{(0)}$ $C_{dB}^{(0)}$ $(\Delta^d C_{dG})$ $(\Delta^d C_{dW})$ $(\Delta^d C_{dB})$	cuW0 cdB0 DeltadcdG DeltadcdW DeltadcdB		
	$\mathcal{L}_6^{(7)}$	$C_{Hl}^{(1)}$ $C_{Hl}^{(3)}$ $(\Delta^u C_{Hq}^{(1)})$ (ΔC_{Hu})	ch11 ch13 DeltaucHq1 DeltacHu	$C_{Hq}^{(1)(0)}$ $C_{Hq}^{(3)(0)}$ $(\Delta^d C_{Hq}^{(1)})$ (ΔC_{Hd})	chq10 chq30 DeltadcHq1 DeltacHd	$C_{Hu}^{(0)}$ $C_{Hd}^{(0)}$ $(\Delta^u C_{Hq}^{(3)})$	chU0 chd0 DeltaucHq3	C_{He} $C_{Hud}^{(0)}$ $(\Delta^d C_{Hq}^{(3)})$	che chud0 DeltadcHq3	
		$\mathcal{L}_6^{(8a)}$	C_{ll} C'_{ll} $(\Delta^u C_{qq}^{(1)})$ $(\Delta^u C_{qq}^{(1)'})$ $(\Delta^u C_{lq}^{(1)})$	c11 c111 Deltaucqq1 Deltaucqq11 Deltauclq1	$C_{qq}^{(1)(0)}$ $C_{qq}^{(1)'(0)}$ $(\Delta^d C_{qq}^{(1)})$ $(\Delta^d C_{qq}^{(1)'})$ $(\Delta^d C_{lq}^{(1)})$	cqq10 cqq110 Deltadcqq1 Deltadcqq11 Deltadclq1	$C_{qq}^{(3)(0)}$ $C_{qq}^{(3)'(0)}$ $(\Delta^u C_{qq}^{(3)})$ $(\Delta^u C_{qq}^{(3)'})$ $(\Delta^u C_{lq}^{(3)})$	cqq30 cqq310 Deltaucqq3 Deltaucqq31 Deltauclq3	$C_{lq}^{(1)(0)}$ $C_{lq}^{(3)(0)}$ $(\Delta^d C_{qq}^{(3)})$ $(\Delta^d C_{qq}^{(3)'})$ $(\Delta^d C_{lq}^{(3)})$	clq10 clq30 Deltadcqq3 Deltadcqq31 Deltadclq3
			$\mathcal{L}_6^{(8b)}$	$C_{uu}^{(0)}$ $C'_{uu}^{(0)}$ C_{ee} $(\Delta^u C_{uu})$ $(\Delta^u C_{ud}^{(1)})$	cuu0 cuu10 cee Deltaucuu Deltaucud1	$C_{dd}^{(0)}$ $C'_{dd}^{(0)}$ (ΔC_{eu}) $(\Delta^d C_{dd})$ $(\Delta^d C_{ud}^{(1)})$	cdd0 cdd10 Deltaceu Deltadccd Deltadcud1	$C_{eu}^{(0)}$ $C_{ed}^{(0)}$ (ΔC_{ed}) $(\Delta^u C'_{uu})$ $(\Delta^u C_{ud}^{(8)})$	ceu0 ced0 Deltaced Deltaucuu1 Deltaucud8	$C_{ud}^{(1)(0)}$ $C_{ud}^{(8)(0)}$ $(\Delta^d C'_{dd})$ $(\Delta^d C_{ud}^{(8)})$
$\mathcal{L}_6^{(8c)}$				C_{le} $C_{qu}^{(1)(0)}$ $C_{qu}^{(1)'(0)}$ $(\Delta_1^u C_{qu}^{(1)})$ $(\Delta_1^u C_{qu}^{(8)})$ $(\Delta_1^u C_{qd}^{(1)})$ $(\Delta_1^u C_{qd}^{(8)})$	cle cqu10 cqu110 Delta1ucqu1 Delta1ucqu8 Delta1ucqd1 Delta1ucqd8	$C_{lu}^{(0)}$ $C_{qd}^{(1)(0)}$ $C_{qd}^{(1)'(0)}$ $(\Delta_1^d C_{qu}^{(1)})$ $(\Delta_1^d C_{qu}^{(8)})$ $(\Delta_1^d C_{qd}^{(1)})$ $(\Delta_1^d C_{qd}^{(8)})$	clu0 cqd10 cqd110 Delta1dcqu1 Delta1dcqu8 Delta1dcqd1 Delta1dcqd8	$C_{ld}^{(0)}$ $C_{qu}^{(8)(0)}$ $C_{qu}^{(8)'(0)}$ $(\Delta_2 C_{qu}^{(1)})$ $(\Delta_2 C_{qu}^{(8)})$ $(\Delta_2 C_{qd}^{(1)})$ $(\Delta_2 C_{qd}^{(8)})$	cld0 cqu80 cqu810 Delta2cqu1 Delta2cqu8 Delta2cqd1 Delta2cqd8	$C_{qe}^{(0)}$ $C_{qd}^{(8)(0)}$ $C_{qd}^{(8)'(0)}$ (ΔC_{lu}) (ΔC_{ld}) $(\Delta^u C_{qe})$ $(\Delta^d C_{qe})$
	$\mathcal{L}_6^{(8d)}$			$C_{lequ}^{(1)(0)}$ $C_{quqd}^{(1)}$ $(\Delta^u C_{ledq})$ $(\Delta^u C_{lequ}^{(1)})$	clequ10 cquqd1 Deltaucledq Deltauclequ1	$C_{lequ}^{(3)(0)}$ $C_{quqd}^{(8)}$ $(\Delta^d C_{ledq})$ $(\Delta^d C_{lequ}^{(1)})$	clequ30 cquqd8 Deltadcledq Deltadclequ1	$C_{ledq}^{(0)}$ $C_{quqd}^{(1)'}$ $(\Delta^u C_{lequ}^{(3)})$	cledq0 cquqd11 Deltauclequ3	$C_{quqd}^{(8)'}$ $(\Delta^d C_{lequ}^{(3)})$

Table 21. The 121 independent parameters in $\mathcal{L}_6^{(5,6,7,8)}$ defined in the MFV model, see section 3.3.

Fermionic SMEFT parameters in SMEFTsim top								
$\mathcal{L}_6^{(5)}$	$\text{Re}(C_{eH})_{pp}$	ceHRe [pp]	$\text{Im}(C_{eH})_{pp}$	ceHIm [pp]	$\text{Re} C_{tH}$	ctHRe	$\text{Im} C_{tH}$	ctHIm
	$\text{Re} C_{uH}$	cuHRe	$\text{Im} C_{uH}$	cuHIm	$\text{Re} C_{bH}$	cbHRe	$\text{Im} C_{bH}$	cbHIm
	$\text{Re} C_{dH}$	cdHRe	$\text{Im} C_{dH}$	cdHIm				
$\mathcal{L}_6^{(6)}$	$\text{Re}(C_{eW})_{pp}$	ceWRe [pp]	$\text{Im}(C_{eW})_{pp}$	ceWIm [pp]	$\text{Re}(C_{eB})_{pp}$	ceBRe [pp]	$\text{Im}(C_{eB})_{pp}$	ceBIm [pp]
	$\text{Re} C_{uG}$	cuGRe	$\text{Im} C_{uG}$	cuGIm	$\text{Re} C_{tG}$	ctGRe	$\text{Im} C_{tG}$	ctGIm
	$\text{Re} C_{dG}$	cdGRe	$\text{Im} C_{dG}$	cdGIm	$\text{Re} C_{bG}$	cbGRe	$\text{Im} C_{bG}$	cbGIm
	$\text{Re} C_{uW}$	cuWRe	$\text{Im} C_{uW}$	cuWIm	$\text{Re} C_{tW}$	ctWRe	$\text{Im} C_{tW}$	ctWIm
	$\text{Re} C_{dW}$	cdWRe	$\text{Im} C_{dW}$	cdWIm	$\text{Re} C_{bW}$	cbWRe	$\text{Im} C_{bW}$	cbWIm
	$\text{Re} C_{uB}$	cuBRE	$\text{Im} C_{uB}$	cuBIm	$\text{Re} C_{tB}$	ctBRE	$\text{Im} C_{tB}$	ctBIm
	$\text{Re} C_{dB}$	cdBRE	$\text{Im} C_{dB}$	cdBIm	$\text{Re} C_{bB}$	cbBRE	$\text{Im} C_{bB}$	cbBIm
$\mathcal{L}_6^{(7)}$	$(C_{Hl}^{(1)})_{pp}$	cHl1 [pp]	$(C_{Hl}^{(3)})_{pp}$	cHl3 [pp]	$(C_{He})_{pp}$	cHe [pp]		
	$C_{Hq}^{(1)}$	cHj1	$C_{Hq}^{(1)}$	cHQ1	C_{Hu}	cHu	C_{Ht}	cHt
	$C_{Hq}^{(3)}$	cHj3	$C_{Hq}^{(3)}$	cHQ3	C_{Hd}	cHd	C_{Hb}	cHbq
	$\text{Re} C_{Hud}$	cHudRe	$\text{Im} C_{Hud}$	cHudIm	$\text{Re} C_{Htb}$	cHtbRe	$\text{Im} C_{Htb}$	cHtbIm
$\mathcal{L}_6^{(8a)}$	$(C_l)_{prst}$	c11 [prst]	prst = {1111, 2222, 3333, 1122, 1133, 2233, 1221, 1331, 2332}					
	$(C_{lq}^{(1)})_{pp}$	c1j1 [pp]	$(C_{lq}^{(1)})_{pp}$	cQ11 [pp]	$(C_{lq}^{(3)})_{pp}$	c1j3 [pp]	$(C_{lq}^{(3)})_{pp}$	cQ13 [pp]
	$C_{qq}^{(1,1)}$	cjj11	$C_{qq}^{(1,8)}$	cjj18	$C_{qq}^{(1,1)}$	cQj11	$C_{qq}^{(1,8)}$	cQj18
	$C_{qq}^{(3,1)}$	cjj31	$C_{qq}^{(3,8)}$	cjj38	$C_{qq}^{(3,1)}$	cQj31	$C_{qq}^{(3,8)}$	cQj38
	$C_{QQ}^{(1)}$	cQQ1	$C_{QQ}^{(8)}$	cQQ8				
$\mathcal{L}_6^{(8b)}$	$(C_{ee})_{prst}$	cee [prst]	prst = {1111, 2222, 3333, 1122, 1133, 2233}					
	$(C_{eu})_{pp}$	ceu [pp]	$(C_{et})_{pp}$	cte [pp]	$(C_{ed})_{pp}$	ced [pp]	$(C_{eb})_{pp}$	cbe [pp]
	$C_{uu}^{(1)}$	cuu1	$C_{uu}^{(8)}$	cuu8	$C_{dd}^{(1)}$	cdd1	$C_{dd}^{(8)}$	cdd8
	$C_{tu}^{(1)}$	ctu1	$C_{tu}^{(8)}$	ctu8	$C_{bd}^{(1)}$	cbd1	$C_{bd}^{(8)}$	cbd8
	C_{tt}	ctt			C_{bb}	cbb		
	$C_{ud}^{(1)}$	cud1	$C_{td}^{(1)}$	ctd1	$C_{ub}^{(1)}$	cbu1	$C_{tb}^{(1)}$	ctb1
	$C_{ud}^{(8)}$	cud8	$C_{td}^{(8)}$	ctd8	$C_{ub}^{(8)}$	cbu8	$C_{tb}^{(8)}$	ctb8
	$\text{Re}(C_{utbd}^{(1)})$	cutbd1Re	$\text{Im}(C_{utbd}^{(1)})$	cutbd1Im	$\text{Re}(C_{utbd}^{(8)})$	cutbd8Re	$\text{Im}(C_{utbd}^{(8)})$	cutbd8Im
$\mathcal{L}_6^{(8c)}$	$(C_{le})_{prst}$	c1e [prst]	prst = {1111, 2222, 3333, 1122, 1133, 2233, 2211, 3311, 3322, 1221, 1331, 2332}					
	$(C_{lu})_{pp}$	clu [pp]	$(C_{lt})_{pp}$	ctl [pp]	$(C_{ld})_{pp}$	cl1 [pp]	$(C_{lb})_{pp}$	clb [pp]
	$(C_{qe})_{pp}$	cje [pp]	$(C_{qe})_{pp}$	cqe [pp]				
	$C_{qu}^{(1)}$	cju1	$C_{qu}^{(1)}$	cQu1	$C_{qt}^{(1)}$	ctj1	$C_{Qt}^{(1)}$	cQt1
	$C_{qu}^{(8)}$	cju8	$C_{qu}^{(8)}$	cQu8	$C_{qt}^{(8)}$	ctj8	$C_{Qt}^{(8)}$	cQt8
	$\text{Re}(C_{qqtu}^{(1)})$	cjqtu1Re	$\text{Im}(C_{qqtu}^{(1)})$	cjqtu1Im	$\text{Re}(C_{qqtu}^{(8)})$	cjqtu8Re	$\text{Im}(C_{qqtu}^{(8)})$	cjqtu8Im
	$C_{qd}^{(1)}$	cjd1	$C_{qd}^{(1)}$	cQd1	$C_{qb}^{(1)}$	cbj1	$C_{Qb}^{(1)}$	cQb1
	$C_{qd}^{(8)}$	cjd8	$C_{qd}^{(8)}$	cQd8	$C_{qb}^{(8)}$	cbj8	$C_{Qb}^{(8)}$	cQb8
$\text{Re}(C_{qQbd}^{(1)})$	cjQbd1Re	$\text{Im}(C_{qQbd}^{(1)})$	cjQbd1Im	$\text{Re}(C_{qQbd}^{(8)})$	cjQbd8Re	$\text{Im}(C_{qQbd}^{(8)})$	cjQbd8Im	
$\mathcal{L}_6^{(8d)}$	$\text{Re}(C_{ledq})_{pp}$	c1edjRe [pp]	$\text{Im}(C_{ledq})_{pp}$	c1edqIm [pp]	$\text{Re}(C_{lebQ})_{pp}$	c1ebQRe [pp]	$\text{Im}(C_{lebQ})_{pp}$	c1ebQIm [pp]
	$\text{Re}(C_{lequ}^{(1)})_{pp}$	c1eju1Re [pp]	$\text{Im}(C_{lequ}^{(1)})_{pp}$	c1eju1Im [pp]	$\text{Re}(C_{leQt}^{(1)})_{pp}$	c1eQt1Re [pp]	$\text{Im}(C_{leQt}^{(1)})_{pp}$	c1eQt1Im [pp]
	$\text{Re}(C_{lequ}^{(3)})_{pp}$	c1eju3Re [pp]	$\text{Im}(C_{lequ}^{(3)})_{pp}$	c1eju3Im [pp]	$\text{Re}(C_{leQt}^{(3)})_{pp}$	c1eQt3Re [pp]	$\text{Im}(C_{leQt}^{(3)})_{pp}$	c1eQt3Im [pp]
	$\text{Re} C_{quqd}^{(1)}$	cjujd1Re	$\text{Im} C_{quqd}^{(1)}$	cjujd1Im	$\text{Re} C_{QtQb}^{(1)}$	cQtQb1Re	$\text{Im} C_{QtQb}^{(1)}$	cQtQb1Im
	$\text{Re} C_{quqd}^{(8)}$	cjujd8Re	$\text{Im} C_{quqd}^{(8)}$	cjujd8Im	$\text{Re} C_{QtQb}^{(8)}$	cQtQb8Re	$\text{Im} C_{QtQb}^{(8)}$	cQtQb8Im
	$\text{Re} C_{quqd}^{(1)'} $	cjujd11Re	$\text{Im} C_{quqd}^{(1)'}$	cjujd11Im	$\text{Re} C_{quqd}^{(8)'}$	cjujd81Re	$\text{Im} C_{quqd}^{(8)'}$	cjujd81Im
	$\text{Re} C_{Qtqd}^{(1)}$	cQtjd1Re	$\text{Im} C_{Qtqd}^{(1)}$	cQtjd1Im	$\text{Re} C_{quQb}^{(1)}$	cjuQb1Re	$\text{Im} C_{quQb}^{(1)}$	cjuQb1Im
	$\text{Re} C_{Qtqd}^{(8)}$	cQtjd8Re	$\text{Im} C_{Qtqd}^{(8)}$	cQtjd8Im	$\text{Re} C_{quQb}^{(8)}$	cjuQb8Re	$\text{Im} C_{quQb}^{(8)}$	cjuQb8Im
	$\text{Re} C_{qtQd}^{(1)}$	cjtQd1Re	$\text{Im} C_{qtQd}^{(1)}$	cjtQd1Im	$\text{Re} C_{Quqb}^{(1)}$	cQujb1Re	$\text{Im} C_{Quqb}^{(1)}$	cQujb1Im
	$\text{Re} C_{qtQd}^{(8)}$	cjtQd8Re	$\text{Im} C_{qtQd}^{(8)}$	cjtQd8Im	$\text{Re} C_{Quqb}^{(8)}$	cQujb8Re	$\text{Im} C_{Quqb}^{(8)}$	cQujb8Im

Table 22. The 260 independent parameters in $\mathcal{L}_6^{(5,6,7,8)}$ defined in the top model, see section 3.4. Lepton flavor indices [pp] always run over {11, 22, 33}. Indices [prst] take the values indicated in line.

Fermionic SMEFT parameters in SMEFTsim topU31								
$\mathcal{L}_6^{(5)}$	Re C_{eH}	ceHRe	Im C_{eH}	ceHIm	Re C_{tH}	ctHRe	Im C_{tH}	ctHIm
	Re C_{uH}	cuHRe	Im C_{uH}	cuHIm	Re C_{bH}	cbHRe	Im C_{bH}	cbHIm
	Re C_{dH}	cdHRe	Im C_{dH}	cdHIm				
$\mathcal{L}_6^{(6)}$	Re C_{eW}	ceWRe	Im C_{eW}	ceWIm	Re C_{eB}	ceBRe	Im C_{eB}	ceBIm
	Re C_{uG}	cuGRe	Im C_{uG}	cuGIm	Re C_{tG}	ctGRe	Im C_{tG}	ctGIm
	Re C_{dG}	cdGRe	Im C_{dG}	cdGIm	Re C_{bG}	cbGRe	Im C_{bG}	cbGIm
	Re C_{uW}	cuWRe	Im C_{uW}	cuWIm	Re C_{tW}	ctWRe	Im C_{tW}	ctWIm
	Re C_{dW}	cdWRe	Im C_{dW}	cdWIm	Re C_{bW}	cbWRe	Im C_{bW}	cbWIm
	Re C_{uB}	cuBRe	Im C_{uB}	cuBIm	Re C_{tB}	ctBRe	Im C_{tB}	ctBIm
	Re C_{dB}	cdBRe	Im C_{dB}	cdBIm	Re C_{bB}	cbBRe	Im C_{bB}	cbBIm
$\mathcal{L}_6^{(7)}$	$C_{Hl}^{(1)}$	cHl1	$C_{Hl}^{(3)}$	cHl3	C_{He}	cHe		
	$C_{Hq}^{(1)}$	cHj1	$C_{HQ}^{(1)}$	cHQ1	C_{Hu}	CHu	C_{Ht}	cHt
	$C_{Hq}^{(3)}$	cHj3	$C_{HQ}^{(3)}$	cHQ3	C_{Hd}	CHd	C_{Hb}	CHbq
	Re C_{Hud}	CHudRe	Im C_{Hud}	CHudIm	Re C_{Htb}	CHtbRe	Im C_{Htb}	CHtbIm
$\mathcal{L}_6^{(8a)}$	C_{ll}	c11	C'_{ll}	c111	$C_{QQ}^{(1)}$	cQQ1	$C_{QQ}^{(8)}$	cQQ8
	$C_{lq}^{(1)}$	clj1	$C_{lQ}^{(1)}$	cQ11	$C_{lq}^{(3)}$	clj3	$C_{lQ}^{(3)}$	cQ13
	$C_{qq}^{(1,1)}$	cjj11	$C_{qq}^{(1,8)}$	cjj18	$C_{Qq}^{(1,1)}$	cQj11	$C_{Qq}^{(1,8)}$	cQj18
	$C_{qq}^{(3,1)}$	cjj31	$C_{qq}^{(3,8)}$	cjj38	$C_{Qq}^{(3,1)}$	cQj31	$C_{Qq}^{(3,8)}$	cQj38
$\mathcal{L}_6^{(8b)}$	C_{eu}	ceu	C_{et}	cte	C_{ed}	ced	C_{eb}	cbe
	$C_{uu}^{(1)}$	cuu1	$C_{uu}^{(8)}$	cuu8	$C_{dd}^{(1)}$	cdd1	$C_{dd}^{(8)}$	cdd8
	$C_{tu}^{(1)}$	ctu1	$C_{tu}^{(8)}$	ctu8	$C_{bd}^{(1)}$	cbd1	$C_{bd}^{(8)}$	cbd8
	C_{tt}	ctt	C_{bb}	cbb	C_{ee}	cee		
	$C_{ud}^{(1)}$	cud1	$C_{td}^{(1)}$	ctd1	$C_{ub}^{(1)}$	cbu1	$C_{tb}^{(1)}$	ctb1
	$C_{ud}^{(8)}$	cud8	$C_{td}^{(8)}$	ctd8	$C_{ub}^{(8)}$	cbu8	$C_{tb}^{(8)}$	ctb8
	Re($C_{qQbd}^{(1)}$)	cjQbd1Re	Im($C_{qQbd}^{(1)}$)	cjQbd1Im	Re($C_{qQbd}^{(8)}$)	cjQbd8Re	Im($C_{qQbd}^{(8)}$)	cjQbd8Im
$\mathcal{L}_6^{(8c)}$	C_{le}	c1e	C_{qe}	cje	C_{Qe}	cQe		
	C_{lu}	clu	C_{lt}	ctl	C_{ld}	c1d	C_{lb}	cb1
	$C_{qu}^{(1)}$	cju1	$C_{Qu}^{(1)}$	cQu1	$C_{qt}^{(1)}$	ctj1	$C_{Qt}^{(1)}$	cQt1
	$C_{qu}^{(8)}$	cju8	$C_{Qu}^{(8)}$	cQu8	$C_{qt}^{(8)}$	ctj8	$C_{Qt}^{(8)}$	cQt8
	Re($C_{qQtu}^{(1)}$)	cjQtu1Re	Im($C_{qQtu}^{(1)}$)	cjQtu1Im	Re($C_{qQtu}^{(8)}$)	cjQtu8Re	Im($C_{qQtu}^{(8)}$)	cjQtu8Im
	$C_{qd}^{(1)}$	cjd1	$C_{Qd}^{(1)}$	cQd1	$C_{qb}^{(1)}$	cbj1	$C_{Qb}^{(1)}$	cQb1
	$C_{qd}^{(8)}$	cjd8	$C_{Qd}^{(8)}$	cQd8	$C_{qb}^{(8)}$	cbj8	$C_{Qb}^{(8)}$	cQb8
Re($C_{qQbd}^{(1)}$)	cjQbd1Re	Im($C_{qQbd}^{(1)}$)	cjQbd1Im	Re($C_{qQbd}^{(8)}$)	cjQbd8Re	Im($C_{qQbd}^{(8)}$)	cjQbd8Im	
$\mathcal{L}_6^{(8d)}$	Re C_{ledq}	cledjRe	Im C_{ledq}	cledqIm	Re C_{lebQ}	clebQRe	Im C_{lebQ}	clebQIm
	Re $C_{lequ}^{(1)}$	c1eju1Re	Im $C_{lequ}^{(1)}$	c1eju1Im	Re $C_{leQt}^{(1)}$	c1eQt1Re	Im $C_{leQt}^{(1)}$	c1eQt1Im
	Re $C_{lequ}^{(3)}$	c1eju3Re	Im $C_{lequ}^{(3)}$	c1eju3Im	Re $C_{leQt}^{(3)}$	c1eQt3Re	Im $C_{leQt}^{(3)}$	c1eQt3Im
	Re $C_{quqd}^{(1)}$	cjujd1Re	Im $C_{quqd}^{(1)}$	cjujd1Im	Re $C_{QtQb}^{(1)}$	cQtQb1Re	Im $C_{QtQb}^{(1)}$	cQtQb1Im
	Re $C_{quqd}^{(8)}$	cjujd8Re	Im $C_{quqd}^{(8)}$	cjujd8Im	Re $C_{QtQb}^{(8)}$	cQtQb8Re	Im $C_{QtQb}^{(8)}$	cQtQb8Im
	Re $C_{quqd}^{(1)'} $	cjujd11Re	Im $C_{quqd}^{(1)'}$	cjujd11Im	Re $C_{quqd}^{(8)'}$	cjujd81Re	Im $C_{quqd}^{(8)'}$	cjujd81Im
	Re $C_{Qtqd}^{(1)}$	cQtjd1Re	Im $C_{Qtqd}^{(1)}$	cQtjd1Im	Re $C_{quQb}^{(1)}$	cjuQb1Re	Im $C_{quQb}^{(1)}$	cjuQb1Im
	Re $C_{Qtqd}^{(8)}$	cQtjd8Re	Im $C_{Qtqd}^{(8)}$	cQtjd8Im	Re $C_{quQb}^{(8)}$	cjuQb8Re	Im $C_{quQb}^{(8)}$	cjuQb8Im
	Re $C_{qtQd}^{(1)}$	cjtQd1Re	Im $C_{qtQd}^{(1)}$	cjtQd1Im	Re $C_{Quqb}^{(1)}$	cQujb1Re	Im $C_{Quqb}^{(1)}$	cQujb1Im
	Re $C_{qtQd}^{(8)}$	cjtQd8Re	Im $C_{qtQd}^{(8)}$	cjtQd8Im	Re $C_{Quqb}^{(8)}$	cQujb8Re	Im $C_{Quqb}^{(8)}$	cQujb8Im

Table 23. The 167 independent parameters in $\mathcal{L}_6^{(5,6,7,8)}$ defined in the topU31 model, see section 3.4.

E Comparison to other SMEFT UFO models

In this section we compare `SMEFTsim` with other UFO models dedicated to SMEFT studies, and provide a mapping of the common parameters. For the time being, the comparison is restricted to `dim6top` [36, 121] and `SMEFT@NLO` [37, 122], that are both based on the Warsaw basis.

For each model we summarize the main features and provide conversion tables with the parameters defined in `SMEFTsim`. To our knowledge, `SMEFTsim` is currently the only publicly available UFO model that implements linearized SMEFT corrections to propagators.

E.1 `dim6top`

`dim6top` [36, 121] contains LO UFO models dedicated to EFT studies in the top sector. Here we refer specifically to `dim6top_LO_UFO` and `dim6top_LO_UFO_each_coupling_order` in the version published in May 2020.

Flavor structure. `dim6top` is based on the recommendations provided in ref. [36], and it assumes a $U(2)^3$ flavor symmetry in the quark sector and a $(U(1)_{l+e})^3$ in the lepton sector. $U(2)^3$ breaking terms are also available and they are implemented explicitly, i.e. without promoting the quark Yukawas to spurions of the flavor symmetry. Contractions inducing both flavor-conserving and violating neutral currents are included.

All fermion masses and Yukawa couplings are neglected, except those of the top and bottom quarks of the tau lepton. The CKM is taken to be the unit matrix.

Operators implemented. `dim6top` contains only operators that modify the interactions of the top quark, and CP violating terms are included. Most operator definitions are identical to those in `SMEFTsim top`, `topU31`. In a few cases, the invariants implemented differ by a Fierz rotation, as detailed in ref. [36].

Input parameters. Both input schemes $\{\alpha_{em}, m_Z, G_F\}$ and $\{m_W, m_Z, G_F\}$ are supported in `dim6top`. Since purely bosonic and leptonic operators are omitted, this only affects the numerical values assigned to the SM parameters and not the dependence on the Wilson coefficients.

SM loop-generated Higgs couplings. Not implemented.

`dim6top` matches very closely the `top` and `topU31` versions of `SMEFTsim` and it can also be mapped to the `general` one. A correspondence with other flavor versions of `SMEFTsim` can only be established partially, due to incompatibilities in the assumed flavor structure.

The mapping between Wilson coefficients defined in `dim6top` and in the `top`, `topU31` versions of `SMEFTsim` is provided in tables 24, 25. The mapping to the `general` version of `SMEFTsim` is provided in tables 26, 27. In both cases, the first table contains parameters with a one-to-one correspondence, while the second contains parameters that require a basis rotation. For example, the point `cQ1M1=1`, `cQ131=3` in `dim6top` corresponds to `cQ1111=4`, `cQ1311=3` (or `clq1Re1133=4`, `clq3Re1133=3`) in `SMEFTsim`.

Overall minus signs in the mapping are due to the fact that `dim6top` and `SMEFTsim` use opposite sign conventions for the definition of covariant derivatives. Although the operator definitions are identical, the *relative* sign between the \mathcal{L}_6 contribution and the corresponding SM coupling is flipped in a few cases. The physics results are identical in both models once this is accounted for. The presence of explicit Yukawa couplings in the tables is due to the different treatment of flavor symmetry breaking terms.

Wilson coefficients inducing flavor-changing neutral currents can be mapped to parameters in `SMEFTsim general`, and the corresponding tables are available upon request.

E.2 SMEFT@NLO

`SMEFT@NLO` [37, 122] is equipped for NLO QCD calculations in `MadGraph5_aMC@NLO`. Here we compare specifically to `SMEFTatNLO v1.0` published in September 2020.

Flavor structure. `SMEFT@NLO` assumes a flavor symmetry $U(2)_q \times U(3)_d \times U(2)_u$ in the quark sector and $U(1)^3$ in the lepton sector, which is the same as in `SMEFTsim top` and in `dim6top`, except for the treatment of down quarks.

All fermion masses and Yukawa couplings are neglected, except those of the top quark.

Operators implemented. `SMEFT@NLO` contains all the operators in classes (1)-(7) and those in class (8) that contain a top quark. Terms that violate the flavor symmetry have been consistently dropped. CP violating terms are omitted.

Input parameters. `SMEFT@NLO` implements the $\{m_W, m_Z, G_F\}$ input scheme.

SM loop-generated Higgs couplings. Higgs couplings in the $m_t \rightarrow \infty$ limit are not implemented, but Higgs-gluon interactions can be fully reproduced at 1-loop in QCD.

Given its flavor structure, `SMEFT@NLO` can be directly mapped to `SMEFTsim` in the `top`, `topU31` and `general` versions. The mapping of Wilson coefficients between `SMEFT@NLO` and the `top`, `topU31` versions of `SMEFTsim` is provided in tables 28, 29. The mapping to the `general` version is provided in tables 30, 31. In both cases, the first table contains the mapping of parameters with a one-to-one correspondence, while the second contains parameters that require a basis rotation.

As for `dim6top`, the sign convention used in `SMEFT@NLO` is the opposite compared to `SMEFTsim`, which leads to some minus signs in the conversion.

SMEFTsim top vs dim6top							
class	SMEFTsim \leftrightarrow dim6top		SMEFTsim \leftrightarrow dim6top		SMEFTsim \leftrightarrow dim6top		
$\mathcal{L}_6^{(5)}$	ctHRe	ctp	ctHIm	ctpI			
$\mathcal{L}_6^{(6)}$	ctGRe	- ctG	ctGIm	- ctGI			
	cbWRe	- cbW	cbWIm	- cbWI			
$\mathcal{L}_6^{(7)}$	cHt	cpt	cHbq	cpb			
	cHtbRe	cptb	cHtbIm	cptbI			
$\mathcal{L}_6^{(8a)}$	cQQ1	$\frac{1}{2}$ cQQ1	cQj11	cQq11	cQj31	cQq13	
	cQQ8	$\frac{1}{2}$ cQQ8	cQj18	cQq81	cQj38	cQq83	
$\mathcal{L}_6^{(8b)}$	ctt	ctt1	ctu1	ctu1	ctu8	ctu8	
	ctd1	ctd1	ctb8	ctb8	ctd8	ctd8	
	ctb1	ctb1	cte[pp]	cte[p]			
$\mathcal{L}_6^{(8c)}$	cQu1	cQu1	ctj1	ctq1	cQt1	cQt1	
	cQu8	cQu8	ctj8	ctq8	cQt8	cQt8	
	cQd1	cQd1	cQb1	cQb1	cQb8	cQb8	
	cQd8	cQd8	cQe[pp]	cQe[p]	ct1[pp]	ct1[p]	
$\mathcal{L}_6^{(8d)}$	clebQRe[pp]	cb1S[p]	c1eQt1Re[pp]	ct1S[p]	c1eQt3Re[pp]	ct1T[p]	
	clebQIm[pp]	cb1SI[p]	c1eQt1Im[pp]	ct1SI[p]	c1eQt3Im[pp]	ct1TI[p]	
	cQtQb1Re	cQtQb1	cQtQb8Re	cQtQb8			
	cQtQb1Im	cQtQb1I	cQtQb8Im	cQtQb8I			
SMEFTsim topU31 vs dim6top							
$\mathcal{L}_6^{(8b)}$	cte	cte[p]					
$\mathcal{L}_6^{(8c)}$	cQe	cQe[p]	ct1	ct1[p]			
$\mathcal{L}_6^{(8d)}$	clebQRe	y1[p] cb1S[p]	c1eQt1Re	y1[p] ct1S[p]	c1eQt3Re	y1[p] ct1T[p]	
	clebQIm	y1[p] cb1SI[p]	c1eQt1Im	y1[p] ct1SI[p]	c1eQt3Im	y1[p] ct1TI[p]	

Table 24. Upper panel: conversion table between the SMEFT parameters defined in the `top` version of SMEFTsim and in `dim6top`: parameters with a one-to-one translation. Lower panel: conversion table between the `topU31` version of SMEFTsim and `dim6top`. Only leptonic coefficients are reported, as the other parameters behave identically to the `top` case. Lepton flavor indices p take values $p = \{1, 2, 3\}$, light quark indices r take values $r = \{1, 2\}$.

SMEFTsim top vs dim6top					
class	SMEFTsim \rightarrow dim6top		dim6top \rightarrow SMEFTsim		
$\mathcal{L}_6^{(6)}$	- ctWRe	ctW	- ctW	ctWRe	
	- ctWIm	ctWI	- ctWI	ctWIm	
	-ctWRe c_θ + ctBRe s_θ	ctZ	ctZ/ s_θ - ctW/ t_θ	ctBRe	
	-ctWIm c_θ + ctBIm s_θ	ctZI	ctZI/ s_θ - ctWI/ t_θ	ctBIm	
$\mathcal{L}_6^{(7)}$	cHQ3	cpQ3	cpQ3 + cpQM	cHQ1	
	cHQ1 - cHQ3	cpQM	cpQ3	cHQ3	
$\mathcal{L}_6^{(8a)}$	cQ11[pp] - cQ13[pp]	cQ1M[p]	cQ13[p] + cQ1M[p]	cQ11[pp]	
	cQ13[pp]	cQ13[p]	cQ13[p]	cQ13[pp]	
$\mathcal{L}_6^{(8b)}$	$\frac{1}{3}$ cutbd1Re + $\frac{4}{9}$ cutbd8Re	yu[r] yd[s] cbtud1	$\frac{1}{3}$ cbtud1 + $\frac{4}{9}$ cbtud8	$\frac{1}{yu[r] yd[s]}$ cutbd1Re	
	$\frac{1}{3}$ cutbd1Im + $\frac{4}{9}$ cutbd8Im	yu[r] yd[s] cbtud1I	$\frac{1}{3}$ cbtud1I + $\frac{4}{9}$ cbtud8I	$\frac{1}{yu[r] yd[s]}$ cutbd1Im	
	2 cutbd1Re - $\frac{1}{3}$ cutbd8Re	yu[r] yd[s] cbtud8	2 cbtud1 - $\frac{1}{3}$ cbtud8	$\frac{1}{yu[r] yd[s]}$ cutbd8Re	
	2 cutbd1Im - $\frac{1}{3}$ cutbd8Im	yu[r] yd[s] cbtud8I	2 cbtud1I - $\frac{1}{3}$ cbtud8I	$\frac{1}{yu[r] yd[s]}$ cutbd8Im	
$\mathcal{L}_6^{(8c)}$	$-\frac{2}{3}$ cjQtu1Re - $\frac{8}{9}$ cjQtu8Re	yu[r] ctQqu1	$-\frac{1}{6}$ ctQqu1 - $\frac{2}{9}$ ctQqu8	$\frac{1}{yu[r]}$ cjQtu1Re	
	$-\frac{2}{3}$ cjQtu1Im - $\frac{8}{9}$ cjQtu8Im	yu[r] ctQqu1I	$-\frac{1}{6}$ ctQqu1I - $\frac{2}{9}$ ctQqu8I	$\frac{1}{yu[r]}$ cjQtu1Im	
	- 4 cjQtu1Re + $\frac{2}{3}$ cjQtu8Re	yu[r] ctQqu8	- ctQqu1 + $\frac{1}{6}$ ctQqu8	$\frac{1}{yu[r]}$ cjQtu8Re	
	- 4 cjQtu1Im + $\frac{2}{3}$ cjQtu8Im	yu[r] ctQqu8I	- ctQqu1I + $\frac{1}{6}$ ctQqu8I	$\frac{1}{yu[r]}$ cjQtu8Im	
	$-\frac{2}{3}$ cjQbd1Re - $\frac{8}{9}$ cjQbd8Re	yd[r] cbQqd1	$-\frac{1}{6}$ cbQqd1 - $\frac{2}{9}$ cbQqd8	$\frac{1}{yd[r]}$ cjQbd1Re	
	$-\frac{2}{3}$ cjQbd1Im - $\frac{8}{9}$ cjQbd8Im	yd[r] cbQqd1I	$-\frac{1}{6}$ cbQqd1I - $\frac{2}{9}$ cbQqd8I	$\frac{1}{yd[r]}$ cjQbd1Im	
	- 4 cjQbd1Re + $\frac{2}{3}$ cjQbd8Re	yd[r] cbQqd8	- cbQqd1 + $\frac{1}{6}$ cbQqd8	$\frac{1}{yd[r]}$ cjQbd8Re	
	- 4 cjQbd1Im + $\frac{2}{3}$ cjQbd8Im	yd[r] cbQqd8I	- cbQqd1I + $\frac{1}{6}$ cbQqd8I	$\frac{1}{yd[r]}$ cjQbd8Im	
$\mathcal{L}_6^{(8d)}$	cQtjd1Re - $\frac{1}{6}$ cjtQd1Re - $\frac{2}{9}$ cjtQd8Re	yd[r] cQtqd1	cQtqd1 - 4 cQtqd1T	$\frac{1}{yd[r]}$ cQtjd1Re	
	cQtjd1Im - $\frac{1}{6}$ cjtQd1Im - $\frac{2}{9}$ cjtQd8Im	yd[r] cQtqd1I	cQtqd1I - 4 cQtqd1TI	$\frac{1}{yd[r]}$ cQtjd1Im	
	cQtjd8Re - cjtQd1Re + $\frac{1}{6}$ cjtQd8Re	yd[r] cQtqd8	cQtqd8 - 4 cQtqd8T	$\frac{1}{yd[r]}$ cQtjd8Re	
	cQtjd8Im - cjtQd1Im + $\frac{1}{6}$ cjtQd8Im	yd[r] cQtqd8I	cQtqd8I - 4 cQtqd8TI	$\frac{1}{yd[r]}$ cQtjd8Im	
	$-\frac{1}{24}$ cjtQd1Re - $\frac{1}{18}$ cjtQd8Re	yd[r] cQtqd1T	$-\frac{8}{3}$ cQtqd1T - $\frac{32}{9}$ cQtqd8T	$\frac{1}{yd[r]}$ cjtQd1Re	
	$-\frac{1}{24}$ cjtQd1Im - $\frac{1}{18}$ cjtQd8Im	yd[r] cQtqd1TI	$-\frac{8}{3}$ cQtqd1TI - $\frac{32}{9}$ cQtqd8TI	$\frac{1}{yd[r]}$ cjtQd1Im	
	$-\frac{1}{4}$ cjtQd1Re + $\frac{1}{24}$ cjtQd8Re	yd[r] cQtqd8T	- 16 cQtqd1T + $\frac{8}{3}$ cQtqd8T	$\frac{1}{yd[r]}$ cjtQd8Re	
	$-\frac{1}{4}$ cjtQd1Im + $\frac{1}{24}$ cjtQd8Im	yd[r] cQtqd8TI	- 16 cQtqd1TI + $\frac{8}{3}$ cQtqd8TI	$\frac{1}{yd[r]}$ cjtQd8Im	
	cjuQb1Re - $\frac{1}{6}$ cQujb1Re - $\frac{2}{9}$ cQujb8Re	yu[r] cQbqu1	cQbqu1 - 4 cQbqu1T	$\frac{1}{yu[r]}$ cjuQb1Re	
	cjuQb1Im - $\frac{1}{6}$ cQujb1Im - $\frac{2}{9}$ cQujb8Im	yu[r] cQbqu8	cQbqu1I - 4 cQbqu1TI	$\frac{1}{yu[r]}$ cjuQb1Im	
	cjuQb8Re - cQujb1Re + $\frac{1}{6}$ cQujb8Re	yu[r] cQbqu1I	cQbqu8 - 4 cQbqu8T	$\frac{1}{yu[r]}$ cjuQb8Re	
	cjuQb8Im - cQujb1Im + $\frac{1}{6}$ cQujb8Im	yu[r] cQbqu8I	cQbqu8I - 4 cQbqu8TI	$\frac{1}{yu[r]}$ cjuQb8Im	
	$-\frac{1}{24}$ cQujb1Re - $\frac{1}{18}$ cQujb8Re	yu[r] cQbqu1T	$-\frac{8}{3}$ cQbqu1T - $\frac{32}{9}$ cQbqu8T	$\frac{1}{yu[r]}$ cQujb1Re	
	$-\frac{1}{24}$ cQujb1Im - $\frac{1}{18}$ cQujb8Im	yu[r] cQbqu1TI	$-\frac{8}{3}$ cQbqu1TI - $\frac{32}{9}$ cQbqu8TI	$\frac{1}{yu[r]}$ cQujb1Im	
	$-\frac{1}{4}$ cQujb1Re + $\frac{1}{24}$ cQujb8Re	yu[r] cQbqu8T	- 16 cQbqu1T + $\frac{8}{3}$ cQbqu8T	$\frac{1}{yu[r]}$ cQujb8Re	
	$-\frac{1}{4}$ cQujb1Im + $\frac{1}{24}$ cQujb8Im	yu[r] cQbqu8TI	- 16 cQbqu1TI + $\frac{8}{3}$ cQbqu8TI	$\frac{1}{yu[r]}$ cQujb8Im	
	SMEFTsim topU31 vs dim6top				
	$\mathcal{L}_6^{(8a)}$	cQ11 - cQ13	cQ1M[p]	cQ13[p] + cQ1M[p]	cQ11
		cQ13	cQ13[p]	cQ13[p]	cQ13

Table 25. Upper panel: conversion table between the SMEFT parameters defined in the top version of SMEFTsim and in dim6top: parameters that require a basis rotation. Lower panel: conversion table between the topU31 version of SMEFTsim and dim6top. Only leptonic coefficients are reported here, as the other parameters behave identically to the top case. Lepton flavor indices take values $p = \{1, 2, 3\}$. In the notation $yd[r]$, $yu[r]$ etc, the flavor index is that of the associated righthanded light field (d for yd and u for yu).

SMEFTsim general vs dim6top						
class	SMEFTsim \leftrightarrow dim6top		SMEFTsim \leftrightarrow dim6top		SMEFTsim \leftrightarrow dim6top	
$\mathcal{L}_6^{(5)}$	cuHRe33	ctp	cuHIm33	ctpI		
$\mathcal{L}_6^{(6)}$	cuGRe33	- ctG	cuGIm33	- ctGI		
	cdWRe33	- cbW	cdWIm33	- cbWI		
$\mathcal{L}_6^{(7)}$	chURe33	cpt	chDRe33	cpb		
	chUdRe33	cptb	chUdIm33	cptbI		
$\mathcal{L}_6^{(8b)}$	ceu[pp]33	cte[p]	cuuRe3333	ctt1	cud1Re33[rr]	ctd1
	cud1Re3333	ctb1	cud8Re3333	ctb8	cud8Re33[rr]	ctd8
$\mathcal{L}_6^{(8c)}$	cqu1Re33[rr]	cQu1	cqu1Re[rr]33	ctq1	cqu1Re3333	cQt1
	cqu8Re33[rr]	cQu8	cqu8Re[rr]33	ctq8	cqu8Re3333	cQt8
	cqd1Re33[rr]	cQd1	cqd1Re3333	cQb1	cqd8Re3333	cQb8
	cqd8Re33[rr]	cQd8	cqeRe33[pp]	cQe[p]	cluRe[pp]33	ctl[p]
$\mathcal{L}_6^{(8d)}$	clEdqRe[pp]33	cb1S[p]	clequ1Re[pp]33	ct1S[p]	clequ3Re[pp]33	ct1T[p]
	clEdqIm[pp]33	cb1SI[p]	clequ1Im[pp]33	ct1SI[p]	clequ3Im[pp]33	ct1TI[p]
	cquqd1Re3333	cQtQb1	cquqd1Im3333	cQtQb1I		
	cquqd8Re3333	cQtQb8	cquqd8Im3333	cQtQb8I		

Table 26. Conversion table between the SMEFT parameters defined in the `general` version of `SMEFTsim` and in `dim6top`: parameters with one-to-one conversion. Lepton flavor indices p take values in $\{1,2,3\}$. Quark flavor indices r take values in $\{1,2\}$.

SMEFTsim general vs dim6top				
class	SMEFTsim \rightarrow dim6top		dim6top \rightarrow SMEFTsim	
$\mathcal{L}_6^{(6)}$	- cuWRe33	ctW	- ctW	cuWRe33
	- cuWIm33	ctWI	- ctWI	cuWIm33
	- cuWRe33 c_θ + cuBRe33 s_θ	ctZ	ctZ/ s_θ - ctW/ t_θ	cuBRe33
	- cuWIm33 c_θ + cuBIm33 s_θ	ctZI	ctZI/ s_θ - ctWI/ t_θ	cuBIm33
$\mathcal{L}_6^{(7)}$	cHq3Re33	cpQ3	cpQ3 + cpQM	cHq1Re33
	cHq1Re33 - cHq3Re33	cpQM	cpQ3	cHq3Re33
$\mathcal{L}_6^{(8a)}$	clq1Re[pp]33 - clq3Re[pp]33	cQ1M[p]	cQ13[p] + cQ1M[p]	clq1Re[pp]33
	clq3[pp]33	cQ13[p]	cQ13[p]	clq3Re[pp]33
	2 cq1Re3333 - $\frac{2}{3}$ cq3Re3333	cQQ1	$\frac{1}{2}$ cQQ1 + $\frac{1}{24}$ cQQ8	cqq1Re3333
	8 cq3Re3333	cQQ8	$\frac{1}{8}$ cQQ8	cqq3Re3333
	$\frac{1}{3}$ cq1Re[r]33[r] + 2 cq1Re[rr]33 + cq3Re[r]33[r]	cQq11	$\frac{1}{2}$ cQq11 - $\frac{1}{12}$ cQq81	cqq1Re[rr]33
	$\frac{1}{3}$ (cq1Re[r]33[r] - cq3Re[r]33[r]) + 2 cq3Re[rr]33	cQq13	$\frac{1}{8}$ cQq81 + $\frac{3}{8}$ cQq83	cqq1Re[r]33[r]
	2 cq1Re[r]33[r] + 6 cq3Re[r]33[r]	cQq81	$\frac{1}{2}$ cQq13 - $\frac{1}{12}$ cQq83	cqq3Re[rr]33
2 (cq1Re[r]33[r] - cq3Re[r]33[r])	cQq83	$\frac{1}{8}$ (cQq81 - cQq83)	cqq3Re[r]33[r]	
$\mathcal{L}_6^{(8b)}$	$\frac{2}{3}$ cuuRe[r]33[r] + 2 cuuRe[rr]33	ctu1	$\frac{1}{2}$ ctu1 - $\frac{1}{12}$ ctu8	cuuRe[rr]33
	4 cuuRe[r]33[r]	ctu8	$\frac{1}{4}$ ctu8	cuuRe[r]33[r]
	$\frac{1}{3}$ cud1Re[r]33[r] + $\frac{4}{9}$ cud8Re[r]33[r]	cbtud1	$\frac{1}{3}$ cbtud1 + $\frac{4}{9}$ cbtud8	cud1Re[r]33[r]
	$\frac{1}{3}$ cud1Im[r]33[r] + $\frac{4}{9}$ cud8Im[r]33[r]	cbtud1I	$\frac{1}{3}$ cbtud1I + $\frac{4}{9}$ cbtud8I	cud1Im[r]33[r]
	2 cud1Re[r]33[r] - $\frac{1}{3}$ cud8Re[r]33[r]	cbtud8	2 cbtud1 - $\frac{1}{3}$ cbtud8	cud8Re[r]33[r]
	2 cud1Im[r]33[r] - $\frac{1}{3}$ cud8Im[r]33[r]	cbtud8I	2 cbtud1I - $\frac{1}{3}$ cbtud8I	cud8Im[r]33[r]
$\mathcal{L}_6^{(8c)}$	- $\frac{2}{3}$ cqu1Re[r]33[r] - $\frac{8}{9}$ cqu8Re[r]33[r]	ctQqu1	- $\frac{1}{6}$ ctQqu1 - $\frac{2}{9}$ ctQqu8	cqu1Re[r]33[r]
	- $\frac{2}{3}$ cqu1Im[r]33[r] - $\frac{8}{9}$ cqu8Im[r]33[r]	ctQqu1I	- $\frac{1}{6}$ ctQqu1I - $\frac{2}{9}$ ctQqu8I	cqu1Im[r]33[r]
	- 4 cqu1Re[r]33[r] + $\frac{2}{3}$ cqu8Re[r]33[r]	ctQqu8	- ctQqu1 + $\frac{1}{6}$ ctQqu8	cqu8Re[r]33[r]
	- 4 cqu1Im[r]33[r] + $\frac{2}{3}$ cqu8Im[r]33[r]	ctQqu8I	- ctQqu1I + $\frac{1}{6}$ ctQqu8I	cqu8Im[r]33[r]
	- $\frac{2}{3}$ cq1Re[r]33[r] - $\frac{8}{9}$ cq8Re[r]33[r]	cbQqd1	- $\frac{1}{6}$ cbQqd1 - $\frac{2}{9}$ cbQqd8	cq1Re[r]33[r]
	- $\frac{2}{3}$ cq1Im[r]33[r] - $\frac{8}{9}$ cq8Im[r]33[r]	cbQqd1I	- $\frac{1}{6}$ cbQqd1I - $\frac{2}{9}$ cbQqd8I	cq1Im[r]33[r]
	- 4 cq1Re[r]33[r] + $\frac{2}{3}$ cq8Re[r]33[r]	cbQqd8	- cbQqd1 + $\frac{1}{6}$ cbQqd8	cq8Re[r]33[r]
	- 4 cq1Im[r]33[r] + $\frac{2}{3}$ cq8Im[r]33[r]	cbQqd8I	- cbQqd1I + $\frac{1}{6}$ cbQqd8I	cq8Im[r]33[r]
	$\mathcal{L}_6^{(8d)}$	cquqd1Re33[rr] - $\frac{1}{6}$ cquqd1Re[r]33[r] - $\frac{2}{9}$ cquqd8Re[r]33[r]	cQtqd1	cQtqd1 - 4 cQtqd1T
cquqd1Im33[rr] - $\frac{1}{6}$ cquqd1Im[r]33[r] - $\frac{2}{9}$ cquqd8Im[r]33[r]		cQtqd1I	cQtqd1I - 4 cQtqd1TI	cquqd1Im33[rr]
cquqd8Re33[rr] - cquqd1Re[r]33[r] + $\frac{1}{6}$ cquqd8Re[r]33[r]		cQtqd8	cQtqd8 - 4 cQtqd8T	cquqd8Re33[rr]
cquqd8Im33[rr] - cquqd1Im[r]33[r] + $\frac{1}{6}$ cquqd8Im[r]33[r]		cQtqd8I	cQtqd8I - 4 cQtqd8TI	cquqd8Im33[rr]
- $\frac{1}{24}$ cquqd1Re[r]33[r] - $\frac{1}{18}$ cquqd8Re[r]33[r]		cQtqd1T	- $\frac{8}{3}$ cQtqd1T - $\frac{32}{9}$ cQtqd8T	cquqd1Re[r]33[r]
- $\frac{1}{24}$ cquqd1Im[r]33[r] - $\frac{1}{18}$ cquqd8Im[r]33[r]		cQtqd1TI	- $\frac{8}{3}$ cQtqd1TI - $\frac{32}{9}$ cQtqd8TI	cquqd1Im[r]33[r]
- $\frac{1}{4}$ cquqd1Re[r]33[r] + $\frac{1}{24}$ cquqd8Re[r]33[r]		cQtqd8T	- 16 cQtqd1T + $\frac{8}{3}$ cQtqd8T	cquqd8Re[r]33[r]
- $\frac{1}{4}$ cquqd1Im[r]33[r] + $\frac{1}{24}$ cquqd8Im[r]33[r]		cQtqd8TI	- 16 cQtqd1TI + $\frac{8}{3}$ cQtqd8TI	cquqd8Im[r]33[r]
cquqd1Re[rr]33 - $\frac{1}{6}$ cquqd1Re3[rr]3 - $\frac{2}{9}$ cquqd8Re3[rr]3		cQbqu1	cQbqu1 - 4 cQbqu1T	cquqd1Re[rr]33
cquqd1Im[rr]33 - $\frac{1}{6}$ cquqd1Im3[rr]3 - $\frac{2}{9}$ cquqd8Im3[rr]3		cQbqu8	cQbqu1I - 4 cQbqu1TI	cquqd1Im[rr]33
cquqd8Re[rr]33 - cquqd1Re3[rr]3 + $\frac{1}{6}$ cquqd8Re3[rr]3		cQbqu1I	cQbqu8 - 4 cQbqu8T	cquqd8Re[rr]33
cquqd8Im[rr]33 - cquqd1Im3[rr]3 + $\frac{1}{6}$ cquqd8Im3[rr]3		cQbqu8I	cQbqu8I - 4 cQbqu8TI	cquqd8Im[rr]33
- $\frac{1}{24}$ cquqd1Re3[rr]3 - $\frac{1}{18}$ cquqd8Re3[rr]3		cQbqu1T	- $\frac{8}{3}$ cQbqu1T - $\frac{32}{9}$ cQbqu8T	cquqd1Re3[rr]3
- $\frac{1}{24}$ cquqd1Im3[rr]3 - $\frac{1}{18}$ cquqd8Im3[rr]3		cQbqu1TI	- $\frac{8}{3}$ cQbqu1TI - $\frac{32}{9}$ cQbqu8TI	cquqd1Im3[rr]3
- $\frac{1}{4}$ cquqd1Re3[rr]3 + $\frac{1}{24}$ cquqd8Re3[rr]3		cQbqu8T	- 16 cQbqu1T + $\frac{8}{3}$ cQbqu8T	cquqd8Re3[rr]3
- $\frac{1}{4}$ cquqd1Im3[rr]3 + $\frac{1}{24}$ cquqd8Im3[rr]3		cQbqu8TI	- 16 cQbqu1TI + $\frac{8}{3}$ cQbqu8TI	cquqd8Im3[rr]3

Table 27. Conversion table between the SMEFT parameters defined in the `general` version of SMEFTsim and in `dim6top`: parameters that require a basis rotation. Lepton flavor indices p take values in $\{1, 2, 3\}$. Quark flavor indices r take values in $\{1, 2\}$.

SMEFTsim top vs SMEFT@NLO						
class	SMEFTsim ↔ SMEFT@NLO		SMEFTsim ↔ SMEFT@NLO		SMEFTsim ↔ SMEFT@NLO	
$\mathcal{L}_6^{(1)}$	cG	- gs cG	cW	- cWWW		
$\mathcal{L}_6^{(2,3)}$	cH	cp	cHbox	cdp	cHDD	cpDC
$\mathcal{L}_6^{(4)}$	cHG	cpG	cHW	cpW	cHB	cpBB
	cHWB	cpWB				
$\mathcal{L}_6^{(5)}$	ctHRe	ctp				
$\mathcal{L}_6^{(6)}$	ctGRe	- gs ctG				
$\mathcal{L}_6^{(7)}$	cHl1[pp]	cpl[p]	cHl3[pp]	c3pl[p]	cHd = cHbq	cpd
	cHe11	cpe	cHe22	cpmu	cHe33	cpta
	cHu	cpu	cHt	cpt		
$\mathcal{L}_6^{(8a)}$	cQj11	cQq11	cQj18	cQq81	cQj31	cQq13
	cQj38	cQq83	cQQ1	$\frac{1}{2}$ cQQ1	cQQ8	$\frac{1}{2}$ cQQ8
	c11[pppp]	c11[pppp]	c11[prrr]	2 c11[prrr]	c11[prrp]	2 c11[prrp]
$\mathcal{L}_6^{(8b)}$	cte[pp]	cte[p]	ctu1	ctu1	ctu8	ctu8
	ctt	ctt1	ctd1 = ctb1	ctd1	ctd8 = ctb8	ctd8
$\mathcal{L}_6^{(8c)}$	ctl[pp]	ctl[p]	cQe[pp]	cQe[p]		
	cQt1	cQt1	cQt8	cQt8		
	cQu1	cQu1	cQd1 = cQb1	cQd1	ctj1	ctq1
	cQu8	cQu8	cQd8 = cQb8	cQd8	ctj8	ctq8
$\mathcal{L}_6^{(8d)}$	c1eQt1Re33	ctlS3	c1eQt3Re33	ctlT3	c1ebQRe33	cb1S3

SMEFTsim topU31 vs SMEFT@NLO						
class	SMEFTsim ↔ SMEFT@NLO		SMEFTsim ↔ SMEFT@NLO		SMEFTsim ↔ SMEFT@NLO	
$\mathcal{L}_6^{(7)}$	cHe	cpe = cpmu = cpta	cHl1	cpl[p]	cHl3	c3pl[p]
$\mathcal{L}_6^{(8a)}$	c11	c11[pppp] = c11[prrr]				
	c111	c11[pppp] = c11[prrp]				
$\mathcal{L}_6^{(8b)}$	cte	cte[p]				
$\mathcal{L}_6^{(8c)}$	cQe	cQe[p]	ctl	ctl[p]		
$\mathcal{L}_6^{(8d)}$	c1eQt1Re33	yl[3] ctlS3	c1eQt3Re33	yl[3] ctlT3	c1ebQRe33	yl[3] cb1S3

Table 28. Upper panel: conversion table between the SMEFT parameters defined in the top version of SMEFTsim and in SMEFT@NLO v1.0, for parameters that have a one-to-one translation. Lower panel: conversion table between the topU31 version of SMEFTsim and SMEFT@NLO. Only leptonic coefficients are reported, as the other parameters behave identically to the top case. The = sign indicates that all parameters need to be fixed to the same value. Lepton flavor indices p take values $p, r = \{1, 2, 3\}$.

SMEFTsim top vs SMEFT@NLO				
class	SMEFTsim \rightarrow SMEFT@NLO		SMEFT@NLO \rightarrow SMEFTsim	
$\mathcal{L}_6^{(6)}$	- ctWRe	ctW	- ctW	ctWRe
	-ctWRe $c_\theta + ctBRe$ s_θ	ctZ	ctZ/ $s_\theta - ctW/t_\theta$	ctBRe
$\mathcal{L}_6^{(7)}$	cHQ3	cpQ3	cpQ3 + cpQM	cHQ1
	cHQ1 - cHQ3	cpQM	cpQ3	cHQ3
	cHj3	cpq3i	cpq3i + cpqMi	cHj1
	cHj1 - cHj3	cpqMi	cpq3i	cHj3
$\mathcal{L}_6^{(8a)}$	cQ11[pp] - cQ13[pp]	cQ1M[p]	cQ13[p] + cQ1M[p]	cQ11[pp]
	cQ13[pp]	cQ13[p]	cQ13[p]	cQ13[pp]
SMEFTsim topU31 vs SMEFT@NLO				
$\mathcal{L}_6^{(8a)}$	cQ11 - cQ13	cQ1M[p]	cQ13[p] + cQ1M[p]	cQ11
	cQ13	cQ13[p]	cQ13[p]	cQ13

Table 29. Upper panel: conversion table between the SMEFT parameters defined in the `top` version of `SMEFTsim` and in `SMEFT@NLO v1.0`: parameters that require a basis rotation. Lower panel: conversion table between the `topU31` version of `SMEFTsim` and `SMEFT@NLO`. Only leptonic coefficients are reported here, as the other parameters behave identically to the `top` case. Lepton flavor indices take values $p = \{1, 2, 3\}$.

SMEFTsim general vs SMEFT@NLO						
class	SMEFTsim \leftrightarrow SMEFT@NLO		SMEFTsim \leftrightarrow SMEFT@NLO		SMEFTsim \leftrightarrow SMEFT@NLO	
$\mathcal{L}_6^{(1)}$	cG	- gs cG	cW	- cWW		
$\mathcal{L}_6^{(2,3)}$	cH	cp	cHbox	cdp	cHDD	cpDC
$\mathcal{L}_6^{(4)}$	cHG	cpG	cHW	cpW	cHB	cpBB
	cHWB	cpWB				
$\mathcal{L}_6^{(5)}$	cuHRe33	ctp				
$\mathcal{L}_6^{(6)}$	cuGRe33	- gs ctG				
$\mathcal{L}_6^{(7)}$	cH11Re [pp]	cp1 [p]	cH13Re [pp]	c3p1 [p]	cHdRe [rr]=cHdRe33	cpd
	cHeRe11	cpe	cHe2Re2	cpmu	cHeRe33	cp _t a
	cHuRe [rr]	cpu	cHuRe33	cpt		
$\mathcal{L}_6^{(8a)}$	c11Re [prst]	c11 [prst]				
$\mathcal{L}_6^{(8b)}$	ceuRe [pp] 33	cte [p]	cuuRe3333	ctt1	cud1Re33 [rr]=cud1Re3333	ctd1
					cud8Re33 [rr]=cud8Re3333	ctd8
$\mathcal{L}_6^{(8c)}$	c1uRe [pp] 33	ct1 [p]	cqeRe33 [pp]	cQe [p]		
	cqu1Re3333	cQt1	cqu8Re3333	cQt8		
	cqu1Re [rr] 33	ctq1	cqu1Re33 [rr]	cQu1	cqd1Re33 [rr]=cqd1Re3333	cQd1
	cqu8Re [rr] 33	ctq8	cqu8Re33 [rr]	cQu8	cqd8Re33 [rr]=cqd8Re3333	cQd8
$\mathcal{L}_6^{(8d)}$	c1equ1Re3333	ct1S3	c1equ3Re3333	ct1T3	c1edqRe3333	cb1S3

Table 30. Conversion table between the SMEFT parameters defined in the general version of SMEFTsim and in SMEFT@NLO v1.0: set of parameters with one-to-one conversion. Lepton flavor indices p take values in {1,2,3}. Quark indices r take values in {1,2}.

SMEFTsim general vs SMEFT@NLO				
class	SMEFTsim \rightarrow SMEFT@NLO		SMEFT@NLO \rightarrow SMEFTsim	
$\mathcal{L}_6^{(6)}$	- cuWRe33	ctW	- ctW	cuWRe33
	- cuWRe33 c_θ + cuBRe33 s_θ	ctZ	ctZ/ s_θ - ctW/ t_θ	cuBRe33
$\mathcal{L}_6^{(7)}$	chq3Re33	cpQ3	cpQ3 + cpQM	chq1Re33
	chq1Re33 - chq3Re33	cpQM	cpQ3	chq3Re33
	chq3Re[rr]	cpq3i	cpq3i + cpqMi	chq1Re[rr]
	chq1Re[rr] - chq3Re[rr]	cpqMi	cpq3i	chq3Re[rr]
$\mathcal{L}_6^{(8a)}$	clq1Re[pp]33 - clq3Re[pp]33	cQ1M[p]	cQ13[p] + cQ1M[p]	clq1Re[pp]33
	clq3[pp]33	cQ13[p]	cQ13[p]	clq3Re[pp]33
	2 cqq1Re3333 - $\frac{2}{3}$ cqq3Re3333	cQQ1	$\frac{1}{2}$ cQQ1 + $\frac{1}{24}$ cQQ8	cqq1Re3333
	8 cqq3Re3333	cQQ8	$\frac{1}{8}$ cQQ8	cqq3Re3333
	$\frac{1}{3}$ cqq1Re[r]33[r] + 2 cqq1Re[rr]33 + cqq3Re[r]33[r]	cQq11	$\frac{1}{2}$ cQq11 - $\frac{1}{12}$ cQq81	cqq1Re[rr]33
	$\frac{1}{3}$ (cqq1Re[r]33[r] - cqq3Re[r]33[r]) + 2 cqq3Re[rr]33	cQq13	$\frac{1}{8}$ cQq81 + $\frac{3}{8}$ cQq83	cqq1Re[r]33[r]
	2 cqq1Re[r]33[r] + 6 cqq3Re[r]33[r]	cQq81	$\frac{1}{2}$ cQq13 - $\frac{1}{12}$ cQq83	cqq3Re[rr]33
2 (cqq1Re[r]33[r] - cqq3Re[r]33[r])	cQq83	$\frac{1}{8}$ (cQq81 - cQq83)	cqq3Re[r]33[r]	
$\mathcal{L}_6^{(8b)}$	$\frac{2}{3}$ cuuRe[r]33[r] + 2 cuuRe[rr]33	ctu1	$\frac{1}{2}$ ctu1 - $\frac{1}{12}$ ctu8	cuuRe[rr]33
	4 cuuRe[r]33[r]	ctu8	$\frac{1}{4}$ ctu8	cuuRe[r]33[r]

Table 31. Conversion table between the SMEFT parameters defined in the **general** version of SMEFTsim and in SMEFT@NLO v1.0: parameters that require a basis rotation. Lepton flavor indices p take values in $\{1, 2, 3\}$. Quark flavor indices r take values in $\{1, 2\}$.

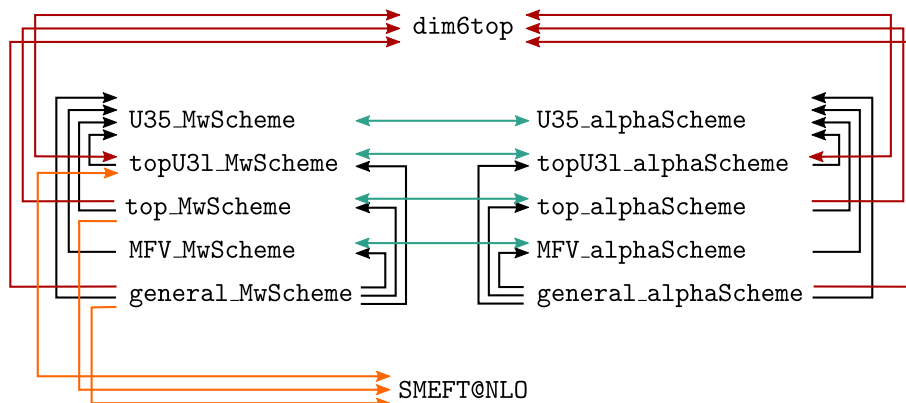


Figure 3. Diagram of the pairwise comparisons performed to validate the UFO models. The arrow indicates the direction in which the SMEFT parameters are mapped.

F Validation of the UFO models

The 10 UFO models contained in the `SMEFTsim` package have been validated following the recommendations of ref. [47]: the procedure relies on pairwise comparisons between models, based on the values returned for a set of squared amplitudes. Each comparison is performed with the dedicated `MadGraph5_aMC@NLO` plugin [123]: given a list of $2 \rightarrow n$ processes and of points in parameter space, the SM squared amplitude $|A_{\text{SM}}|^2$, the pure SM- \mathcal{L}_6 interference $2 \text{Re} A_{\text{SM}} A_6^*$ and the quadratic \mathcal{L}_6 contribution $|A_6|^2$ are calculated at one random phase-space point for each process and parameter point. The validation is considered successful if the squared amplitudes evaluated with each model pair agree within a permille. Larger discrepancies are ignored if they do not show a consistent pattern across different processes and the squared amplitude is $< 10^{-16}$ for both models.

Figure 3 illustrates diagrammatically the set of comparisons performed: the `top`, `topU31` and `general` versions of `SMEFTsim` have been compared to `dim6top` (version of May 2020) and `SMEFT@NLO` (both versions of August 2019 and September 2020, only for models with $\{m_W, m_Z, G_F\}$ scheme). An internal validation was also carried out, comparing models with different flavor assumptions and same input scheme, and vice versa. The arrows in the figure indicate that, in the comparison, the parameters of the first model were mapped onto those of the latter: the flow generally goes towards more restrictive flavor assumptions.

The validation was performed on the processes listed in table 32, that were chosen so as to probe most effective operators independently. All Wilson coefficients have been included in the comparison, with the exception of those inducing flavor-changing neutral currents.

All fermion masses and Yukawa couplings were retained for internal validation, while only those implemented in `dim6top` or `SMEFT@NLO` were included when comparing to these models. CKM mixing has been neglected in all cases.

Linearized propagator corrections have been validated with an analogous procedure, using the processes listed in table 33 and carrying out internal comparisons across models with same inputs and different flavor assumption. Loop-induced SM Higgs couplings have

g g > t t	g g > u u	g g > g g	w+ w- > w+ w-	a w+ > z w+	z z > w+ w-
w+ w- > a a	z z > z z	h h > w+ w-	h h > z z	h h > h h	h w- > t b
h w+ > u d	h w- > mu- vm	h z > b b	h z > s s	h z > t t	h z > u u
h a > b b	h a > t t	h a > d d	h a > c c	h z > mu+ mu-	h z > vt vt
h a > e+ e-	b b > t t	b b > b b	b b > u u	b b > d d	b b > a a
t t > t t	t t > c c	t t > s s	t t > d d	b t > d u	u u > u u
u u > c c	s s > s s	d d > s s	d d > c c	u d > u d	u d > c s
e+ e- > mu+ mu-	ve ve > vm vm	vt vt > vt vt	e+ ta- > e+ ta-	mu+ mu- > ta+ ta-	e+ e- > e+ e-
e+ ve > e+ ve	e+ vm > e+ vm	e+ ve > mu+ vm	b t > e- ve	u d > mu+ vm	c s > ta+ vt
d d > e+ e-	c c > mu+ mu-	b b > mu+ mu-	t t > ta+ ta-	b b > ve ve	s s > vt vt
t t > vm vm	u u > ve ve	w+ w- > h d d	w+ w- > h c c	w+ w- > h t t	w+ w- > h b b
w+ w- > a a a	h w+ > a u d	h w- > a e- ve	h w+ > a t b	h h > h u u	h h > h ta+ ta-
h h > h s s	h h > h t t	h h > h b b	h h > z ve ve	h h > z c c	h h > z t t
h h > z e+ e-	h h > z d d	h h > z b b	h h > w+ e- ve	h h > w- u d	h h > w+ t b
h h > w+ w- a	g g > g g g	h h > h h h	h h > h z z		

Table 32. Set of processes that have been used to validate the UFO models.

w+ w- > w+ w-	z z > w+ w-	w+ w- > a a	h h > w+ w-	h h > z z	h h > h h
h w+ > u d	h w- > mu- vm	h z > s s	h z > t t	h z > mu+ mu-	h z > vt vt
b b > t t	b b > b b	t t > t t	b t > d u	u u > u u	s s > s s
d d > c c	u d > c s	e+ e- > mu+ mu-	ve ve > vm vm	e+ ta- > e+ ta-	e+ e- > e+ e-
e+ ve > e+ ve	e+ vm > e+ vm	e+ ve > mu+ vm	b t > e- ve	c s > ta+ vt	c c > mu+ mu-
b b > mu+ mu-	b b > ve ve	s s > vt vt	u u > ve ve	t g > b w+	w+ w- > h d d
w+ w- > h c c	w+ w- > h b b	w+ w- > a a a	h w+ > a u d	h w- > a e- ve	h w+ > a t b
h h > h ta+ ta-	h h > h b b	h h > z c c	h h > z t t	h h > z b b	h h > w+ e- ve
h h > w- u d	h h > w+ w- a				

Table 33. Set of processes that have been used to validate linearized propagator corrections implemented in the UFO models.

g g > t t	g g > u u	g g > g g	a w+ > z w+	z z > w+ w-	w+ w- > a a
h a > b b	h a > d d	h a > c c	h a > e+ e-	b b > a a	t t > z a
w+ w- > h d d	w+ w- > h t t	w+ w- > h b b	w+ w- > a a a	h h > z c c	h h > z e+ e-
h h > z d d	h h > z b b	g g > g g g	h h > h h h		

Table 34. Set of processes that have been used to validate loop-induced SM Higgs couplings implemented in the UFO models.

been validated comparing the SM squared amplitudes at one phase-space point for each of the processes in table 34 with all models.

The output files generated by the `MadGraph5_aMC@NLO` plugin are available at the github repository. No significant residual differences are present between models. Only one exception was observed: potentially large discrepancies are present between `SMEFTsim` (both `top` and `general`) and `dim6top`, for the Wilson coefficients `cQbqu1T`, `cQbqu8T`, `cQtqd1T`, `cQtqd8T` and the associated imaginary parts. These differences have been already noted in the past and are currently not fully understood.

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