smelli – the SMEFT Likelihood

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I present the Python package smelli that implements a global likelihood function in the space of dimension-six Wilson coefficients in the Standard Model Effective Field Theory (SMEFT). The likelihood includes contributions from a large number of flavor and other precision observables, currently 399 in total.
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### 1. Introduction

The Standard Model (SM) of particle physics is an extremely successful model. However, there are several experimental as well as theoretical indications for new physics (NP) beyond the SM. Whether a given NP scenario describes the experimental data better than the SM can be conveniently quantified by the ratio of the NP likelihood $L_{NP}$ and the SM likelihood $L_{SM}$ or, equivalently, by the difference of the log-likelihoods

$$\Delta \log L = \log L_{NP} - \log L_{SM}. \quad (1)$$

These likelihood functions are constructed from a set of measured observables and take into account uncertainties and correlations from both the measurements and the theoretical predictions.

A set of observables for which certain NP scenarios can describe the experimental data considerably better than the SM have been found e.g. in $B$ meson decays. These so-called $B$ anomalies correspond to deviations from the SM predictions in measurements of neutral current $b \to s\ell\ell$ and charged current $b \to c\ell\nu$ transitions. In particular, deviations have been found in

(i) angular observables of $B \to K^* \mu^+\mu^-$ [1–5],

(ii) branching ratios of $B \to K \mu^+\mu^-$, $B \to K^* \mu^+\mu^-$, and $B_s \to \phi \mu^+\mu^-$ [6–8],

(iii) the lepton flavor universality (LFU) observables $R_{K^{(*)}}$ [9–12], which are $\mu/e$ ratios of $B \to K^{(*)}\ell^+\ell^-$ branching ratios,
(iv) the branching ratio of $B_s \to \mu^+\mu^-$ [13–17].

(v) the LFU observables $R_{D^{(*)}}$ [18–25], which are $\tau/e$ and $\tau/\mu$ ratios of $B \to D^{(*)}\ell\nu$ branching ratios.

While (i) and (ii) could be afflicted by underestimated hadronic uncertainties, the observables in (iii), (iv), and (v) are theoretically clean probes of NP [26–28]. Considering the above $B$-decay observables and parameterizing NP in $b \to s\ell\ell$ and $b \to c\ell\nu$ transitions in terms of Wilson coefficients in the Weak Effective Theory (WET), simple one- and two-parameter scenarios show a sizable $\Delta \log L \sim 20$ (cf. e.g. [29–34]).

These intriguing hints for NP have led to extensive model building. In the process, important insights have been gained:

- The fact that NP above the electroweak (EW) scale has to respect SM gauge invariance leads to important correlations between low-energy observables. For example, explanations of $R_{D^{(*)}}$ in terms of left-handed contributions to $b \to c\tau\nu$ imply also contributions to $b \to s\nu\nu$, which are constrained by $B \to K^{(*)}\nu\bar{\nu}$ [35].

- One-loop contributions can have very important effects. This has been observed in models explaining $R_{D^{(*)}}$ and $R_K^{(*)}$ using mostly 3rd generation couplings. They actually modify $\tau$ and $Z$ decays at one loop, which leads to strong constraints [36]. Another example is provided by models explaining $R_{D^{(*)}}$ using a contribution to semi-tauonic operators, which generate an effect in $b \to s\ell\ell$ at one loop [37, 38].

Essentially every model that explains some of the $B$ anomalies predicts deviations from the SM also in other observables. In many cases, this leads to strong constraints or exclusion of a model. So phenomenological analyses that consider only a small set of observables or neglect one-loop contributions are in many cases not sufficient to show that a given model agrees with experimental data better than the SM. In order to show this, it is in general necessary to

- compute all relevant observables $\hat{O}(\vec{\xi})$ (flavor observables, EW precision observables (EWPO), etc.) in terms of the Lagrangian parameters $\vec{\xi}$ of a NP model,

- take into account loop effects when computing the observables,

- compare the theory predictions to experimental data by constructing the NP likelihood $L_{NP}$.

Performing these steps again and again for each single model one wants to analyze is a tedious task. Fortunately, analyses of NP models can be tremendously simplified by making use of the SM effective field theory (SMEFT) in an intermediate step.

2. The SMEFT Likelihood

Assuming that the scale of NP $\Lambda_{NP}$ is considerably larger than the EW scale and EW symmetry breaking is realized linearly, the NP effects in a given observable can be expressed in terms of the Wilson coefficients $C_i$ of the SMEFT, which are defined by the SMEFT Lagrangian [39, 40]

$$L_{SMEFT} = L_{SM} + \sum_{n>4} \sum_i \frac{C_i}{\Lambda_{NP}^{n-4}} O_i,$$

(2)
where $O_i$ are local SM gauge invariant operators constructed from the SM fields and $n$ is their canonical dimension.

The SMEFT is a powerful tool since it can connect the model building at the high scale $\Lambda_{NP}$ to the phenomenology at lower scales without the need to compute hundreds of observables in each model. A phenomenological analysis can be split into

- a model-dependent part that consists of matching the NP model to the SMEFT at the scale $\Lambda_{NP}$,
- the model-independent phenomenology, which corresponds to
  - running down the Wilson coefficients $\tilde{C}$ from $\Lambda_{NP}$ to the low scale at which the observables are computed,
  - predicting all the relevant observables $\tilde{O}(\tilde{C})$ in terms of the Wilson coefficients $\tilde{C}$,
  - constructing the NP likelihood $L_{NP}(\tilde{O}(\tilde{C}))$ that compares the predictions to experimental measurements,
  - computing $\Delta \log L$ using eq. (1) in order to compare the NP model to the SM.

While it might be preferable to perform the model-dependent matching at one-loop, a large number of important one-loop effects is actually already included by the model-independent renormalization group (RG) running and mixing in the SMEFT.

Using the above procedure, a SMEFT likelihood function $L_{NP}(\tilde{C})$ can tremendously simplify analyses of NP models. Many likelihood functions in the SMEFT have been considered in the literature (see e.g. [41–55]). However, most of them are constructed from observables in one or few specific sectors, like EWPO, Higgs physics, top physics, $B$ physics, or lepton flavor violating observables. But as discussed above, NP models generically predict new effects in several observables of various sectors. Furthermore, SMEFT operators belonging to different sectors mix under renormalization. Consequently, to test a NP model, the sectors should not be considered separately. It is in fact necessary to construct the global SMEFT likelihood, taking into account as many observables from as many sectors as possible.

3. The smelli Python package

In [56], we have started constructing a global SMEFT likelihood that is provided by the Python package smelli (SMEFT likelihood). It is based on

- the Python package flavio [57] that can compute hundreds of flavor and other precision observables in and beyond the SM, while properly accounting for theory uncertainties,
- the Wilson coefficient exchange format (WCxf) [58] that is used to represent and exchange large sets of Wilson coefficients in various EFTs and bases,
- the Python package wilson [59] that performs the RG evolution in the SMEFT and the WET as well as the matching between them.

smelli is built upon these tools and implements a SMEFT likelihood function constructed from currently 399 observables. In particular, it includes
• flavor-changing neutral current $B$ decays,
• lepton flavor universality tests in charged- and neutral-current $B$ and $K$ decays,
• meson-antimeson mixing in the $K$, $B$, and $D$ systems,
• charged lepton flavor violating $B$, tau, and muon decays,
• the anomalous magnetic moments of the electron, muon, and tau,
• $Z$ and $W$ pole EWPO,
• nuclear and neutron beta decays,
• Higgs signal strengths.

Given any combination of SMEFT or WET Wilson coefficients, smelli computes the $\Delta \log L$ for each of the above sectors and then sums all of them to obtain the global $\Delta \log L$.

The full global likelihood is work in progress and the development is open to everyone. The open-source code of smelli is available at https://github.com/smelli/smelli.

3.1 Installation

The requirements for smelli are a working installation of Python version 3.5 or above and the Python package manager pip. If both are present, smelli can be installed from the command line by entering

```
python3 -m pip install smelli --user
```

This will download smelli and all its dependencies from the Python package archive (PyPI) and install it in the user’s home directory without requiring root privileges (due to the option --user).

3.2 Using smelli

Like any Python package, smelli can be used

• as a library imported from other scripts,
• directly in the command line interpreter,
• in an interactive session, e.g. in a Jupyter notebook.

How to use smelli is demonstrated in the following with examples from an interactive Jupyter notebook. This notebook is available at https://github.com/peterstangl/smelli-talk. For further information on the features of smelli, see [56] and the API documentation at https://smelli.github.io.

3.2.1 Instantiating the likelihood

The main functionality of smelli is provided by the GlobalLikelihood class. It is imported by

```
In: from smelli import GlobalLikelihood
```

If the GlobalLikelihood class is instantiated without any argument,

```
In: gl = GlobalLikelihood()
```
the likelihood is defined in the space of SMEFT Wilson coefficients in the **Warsaw** basis (for details on the specifications of the supported EFTs and bases, see the WCxf website at [https://wcxf.github.io/bases.html](https://wcxf.github.io/bases.html)). The EFT and basis of a given `GlobalLikelihood` instance can be accessed via its `eft` and `basis` attributes.

In: `gl.eft, gl.basis`

Out: ('SMEFT', 'Warsaw')

In order to create a likelihood function of Wilson coefficients in the WET, one can provide the `eft` and `basis` arguments on instantiation of a `GlobalLikelihood` instance.

In: `gl_wet = GlobalLikelihood(eft='WET', basis='flavio')`

Out: ('WET', 'flavio')

### 3.2.2 Fixing a point in Wilson coefficient space: 3 equivalent ways

The point in the Wilson coefficient space at which the likelihood should be computed is defined using the `parameter_point` method. This method returns an instance of the `GlobalLikelihoodPoint` class that can be used to compute $\Delta \log L$. The values of the Wilson coefficients can be provided in three equivalent ways:

- A dictionary of Wilson coefficients as well as the scale in GeV at which they are defined can be passed directly as arguments.

  In: `pp = gl.parameter_point({lq3_2223: 1e-9}, scale=1/zero.alt3/zero.alt3)`

- An instance of the Wilson class from the `wilson` package can be passed as a single argument.

  In: `from wilson import Wilson
  w = Wilson({'lq3_2223': 1e-9}, scale=1000, eft='SMEFT', basis='Warsaw')
p = gl.parameter_point(w)`

- A WCxf file, e.g. a file in YAML format named `my_wcxf.yaml` and containing

  
  ```yaml
  eft: SMEFT
  basis: Warsaw
  scale: 1000
  values:
  lq3_2223:
    Re: 1e-9
  
  
  ```

  can be read in by providing the path to the file as argument.

  In: `pp = gl.parameter_point('my_wcxf.yaml')`
### 3.2.3 Computing the likelihood

After the Wilson coefficients have been fixed and an instance of GlobalLikelihoodPoint has been created, it can be used to compute $\Delta \log L$. In smelli, the global $\Delta \log L$ is given in terms of the sum of several individual $\Delta \log L$ that are constructed from subsets of observables. To access all these individual $\Delta \log L$, the method log_likelihood_dict can be used. It returns a dictionary containing the names of the individual likelihoods and the corresponding $\Delta \log L$ values. Using the above defined parameter point, one gets

\[
\begin{align*}
\text{In:} & \quad \text{pp.log_likelihood_dict()} \\
\text{Out:} & \quad \{ \\
& \quad 'fast\_likelihood\_quarks.yaml': 18.063309775625527, \\
& \quad 'fast\_likelihood\_leptons.yaml': -7.954151298861234e-05, \\
& \quad 'likelihood\_ewpt.yaml': 0.0019331634397694586, \\
& \quad 'likelihood\_eeew.yaml': -0.0001731988511934901, \\
& \quad 'likelihood\_lept.yaml': 3.7762380644679183e-07, \\
& \quad 'likelihood\_rd\_rds.yaml': 0.27864506193111893, \\
& \quad 'likelihood\_lfu\_fccc.yaml': 0.000502717997831865, \\
& \quad 'likelihood\_lfu\_fcnc.yaml': 3.0607966063245655, \\
& \quad 'likelihood\_bcpv.yaml': 0.013775072147421241, \\
& \quad 'likelihood\_bqnunu.yaml': -0.119578242544371, \\
& \quad 'likelihood\_lfv.yaml': 0.0, \\
& \quad 'likelihood\_zlfv.yaml': 0.0, \\
& \quad 'likelihood\_higgs.yaml': 2.176258307784451e-05, \\
& \quad 'global': 21.299153554766516
\}\end{align*}
\]

While the global $\Delta \log L$ is provided by log_likelihood_dict, its value can also be directly returned using the log_likelihood_global method.

\[
\begin{align*}
\text{In:} & \quad \text{pp.log_likelihood_global()} \\
\text{Out:} & \quad 21.299153554766516
\end{align*}
\]

Apart from $\Delta \log L$, it is also possible to compute the total $\chi^2_{NP}$, defined by

\[
\chi^2_{NP} = -2 \log L_{NP},
\]

where $L_{NP}$ is normalized such that it is 1 if the central values of the theory predictions are equal to the central values of the measurements for all observables. A dictionary containing the individual values of the total $\chi^2_{NP}$ is returned by the chi2_dict method.

\[
\begin{align*}
\text{In:} & \quad \text{pp.chi2_dict()} \\
\text{Out:} & \quad \{ \\
& \quad 'fast\_likelihood\_quarks.yaml': 160.14558316478963, \\
& \quad 'fast\_likelihood\_leptons.yaml': 23.57908813232271, \\
& \quad 'likelihood\_ewpt.yaml': 35.3618189920579, \\
& \quad 'likelihood\_eeew.yaml': 61.19130715429686, \\
& \quad 'likelihood\_lept.yaml': 1.4486600571844703,
\}\end{align*}
\]
These values are particularly useful for computing p-values from the total $\chi^2_{NP}$ and the number of observations. The latter are returned by the `number_observations_dict` method of the `GlobalLikelihood` instance (which can be conveniently accessed using the `likelihood` attribute of the `GlobalLikelihoodPoint` instance).

In: ```python
pp.likelihood.number_observations_dict()
```

Out: ```python
{'fast_likelihood_quarks.yaml': 144,
 'fast_likelihood_leptons.yaml': 7,
 'likelihood_ewpt.yaml': 30,
 'likelihood_eeww.yaml': 48,
 'likelihood_lept.yaml': 2,
 'likelihood_rd_rds.yaml': 11,
 'likelihood_lfu_fccc.yaml': 63,
 'likelihood_lfu_fcnc.yaml': 21,
 'likelihood_bcpv.yaml': 6,
 'likelihood_bqnuu.yaml': 22,
 'likelihood_lfv.yaml': 41,
 'likelihood_zlfv.yaml': 7,
 'likelihood_higgs.yaml': 67,
 'global': 469}
```

Note that here an “observation” is defined as an individual measurement of an observable. Thus, the number of observations is always greater than or equal to the number of observables.

### 3.2.4 Table of observables

`smelli` provides information on individual observables. In particular, the theoretical and experimental central values and uncertainties as well as the pull compared to the SM or the experimental data can be obtained. All this information is contained in an “observable table” that is returned in the form of a Pandas [60, 61] DataFrame object by the method `obstable`.

In: ```python
df = pp.obstable()
```
In a Jupyter notebook, a Pandas DataFrame is shown as a table.

In: \texttt{df} \\

### Out: \\

| experiment | exp. unc. | theory | th. unc. | pull exp. | pull SM |
|------------|----------|--------|----------|-----------|---------|
| \texttt{a\_mu} | 0.01116592 | 6.31304e-10 | 0.00116592 | 4.25176e-10 | 3.492396 -4.46085e-05 |
| \texttt{Rtaul(B->D^*lnu)} | 0.296146 | 0.015608 | 0.244875 | 0 | 3.306066 -0.389707 |
| (\texttt{<dR/dtheta>(ee->WW), 198.38, 0.8, 1.0}) | 6.535 | 0.236 | 7.236 | 0 | 2.970365 0.0112166 |
| \texttt{BR(W->taunu)} | 0.1138 | 0.0021 | 0.108417 | 0 | 2.563346 -0.00503662 |
| \texttt{epsp/eps} | 0.00166382 | 0.000227703 | -3.12549e-05 | 0.000637111 | 2.505372 0.0147821 |
| ... | ... | ... | ... | ... | ... |
| \texttt{BR(tau->phie)} | 0 | 1.88467e-08 | 0 | 0 | 0 |
| \texttt{BR(tau->phimu)} | 0 | 5.10684e-08 | 0 | 0 | 0 |
| \texttt{BR(Z->emu)} | 0 | 2.33904e-07 | 0 | 0 | 0 |
| \texttt{BR(Z->etau)} | 0 | 2.59807e-06 | 0 | 0 | 0 |
| \texttt{BR(Z->mutau)} | 0 | 2.69574e-06 | 0 | 0 | 0 |

399 rows \times 6 columns

The Pandas DataFrame is a convenient object for tabulated data and provides many useful features. E.g. one can sort the rows by the values of a given column,

In: \texttt{df.sort_values('pull SM', ascending=True)[:5]} \\

### Out: \\

| experiment | exp. unc. | theory | th. unc. | pull exp. | pull SM |
|------------|----------|--------|----------|-----------|---------|
| (\texttt{<dBR/dq2>(Bs->phimumu), 1.0, 6.0}) | 2.55342e-08 | 3.72621e-09 | 4.04247e-08 | 6.44267e-09 | 2.0007 -3.24157 |
| (\texttt{<Rmue>(B0->K*ll), 1.1, 6.0}) | 0.681356 | 0.123108 | 0.746295 | 0 | 0.623038 -2.4685 |
| \texttt{BR(Bs->mumu)} | 2.73001e-09 | 3.80964e-10 | 2.73442e-09 | 1.47033e-09 | 0.0108006 -2.29374 |
| (\texttt{<dBR/dq2>(Bs->phimumu), 15.0, 19.0}) | 4.05106e-08 | 5.09449e-09 | 4.08896e-08 | 4.5361e-09 | 0.0555647 -2.21418 |
| (\texttt{<dBR/dq2>(B0->K*mumu), 15.0, 19.0}) | 4.35409e-08 | 3.61869e-09 | 4.35383e-08 | 6.16124e-09 | 0.000370693 -2.20919 |

or select a specific row by its name.

In: \texttt{df.loc[['Rtaul(B->D^*lnu)']]} \\

### Out: \\

| experiment | exp. unc. | theory | th. unc. | pull exp. | pull SM |
|------------|----------|--------|----------|-----------|---------|
| \texttt{Rtaul(B->D^*lnu)} | 0.296146 | 0.015608 | 0.244875 | 0 | 3.306066 -0.389707 |
### 3.2.5 Plots

Given a likelihood function, one common task is to plot this function in a 2D plane. In order to simplify this, smelli provides a method to compute the plot data for all individual likelihoods. For demonstration, it is convenient to define a `GlobalLikelihood` instance for which the likelihood can be computed much faster than in the default case. This can be achieved by considering only a subset of observables, e.g. only EWPO and the Higgs signal strengths.

```python
In: gl_ewpt_higgs = GlobalLikelihood(include_likelihoods=['likelihood_ewpt.yaml', 'likelihood_higgs.yaml'])
```

The next step is to define a function of the two plot parameters that returns a dictionary of Wilson coefficients. This function defines what is actually plotted. It can be a trivial function that takes two Wilson coefficients as arguments and just returns them, but it can also be a complicated function of two NP model parameters that returns a large set of Wilson coefficients depending on these two parameters. As an example, we will reproduce figure 2 of [62] and plot the likelihood in the space of the $S$ and $T$ parameters. They are proportional to the SMEFT Wilson coefficients $C_{\phi WB}$ and $C_{\phi D}$, and their relations are given by

$$
C_{\phi WB} = \frac{g_L g_Y}{16 \pi v^2} S, \quad C_{\phi D} = -\frac{g_L^2 g_Y^2}{2 \pi (g_L^2 + g_Y^2) v^2} T.
$$

Consequently, plugging in the SM parameters, the function that takes $S$ and $T$ as arguments and returns a dictionary of Wilson coefficients can be defined as follows.

```python
In: def wc_fct(S, T):
    return {
        'phiWB': S * 7.643950529889027e-08,
        'phiD': -T * 2.579372285227687e-07,
    }
```

This function can now be used as the first argument of the `plot_data_2d` method of the `GlobalLikelihood` instance. The second argument is the scale at which the Wilson coefficients are defined, followed by the minimum and maximum values for the x- and y-axis. In the function call below, also two optional arguments are given: the number of steps in each direction (`steps = 10` results in plot data computed on a $10 \times 10$ grid), and the number of CPU threads to be used for the computation.

```python
In: plot_data = gl_ewpt_higgs.plot_data_2d(wc_fct, 91.1876, -0.2, 0.2, -0.1, 0.3, steps=10, threads=8,)
```
The `plot_data_2d` method returns a dictionary with the names of the individual likelihoods as keys and values that are again dictionaries. The keys in these latter dictionaries are \( x, y, \) and \( z \) and the values are arrays. Here, \( x \) and \( y \) correspond to the coordinates in the 2D plane and \( z \) to the values of \( \Delta \chi^2 = -2 \Delta \log L \) at these coordinates. The dictionaries with keys \( x, y, \) and \( z \) are constructed in such a way that they can be directly fed to the contour plotting function of the `flavio` package. The relevant submodules for plotting have to be imported from `flavio` and `matplotlib` [63] (on which the `flavio` plotting functions are based on).

```
In: import flavio.plots as fpl
     import matplotlib.pyplot as plt
```

In order to plot \( \Delta \chi^2 \) contours corresponding to a given pull in units of \( \sigma \), the contour levels can be defined using the `flavio` function `delta_chi2`, which takes the number of \( \sigma \) and the number of degrees of freedom as arguments.

```
In: levels_1sig = [fpl.delta_chi2(1, dof=2)]
levels_123sig = [fpl.delta_chi2(n_sigma, dof=2) for n_sigma in (1,2,3)]
```

The data can now be plotted. The function `fpl.contour` is called three times, once for each of the three different likelihoods: Higgs physics, EWPO, and their combination. Furthermore, horizontal and vertical axes as well as labels are added. A value larger than one for the argument `interpolation_factor` of `fpl.contour` makes the contours appear smooth. However, if the plot data has been computed on a small grid, `interpolation_factor` can obscure the fact that the data might be insufficient for a reasonable plot. In fact, for more reasonable plots, the number of `steps` should be increased to at least 20 (but this of course also increases the computing time). From the data computed above, the plot is then generated by the following code.

```
In: plt.figure(figsize=(5,5))
fpl.contour(**plot_data['likelihood_higgs.yaml'], levels=levels_1sig,
           label=r"Higgs ($1\sigma$)", interpolation_factor=9,
           color='C0')
fpl.contour(**plot_data['likelihood_ewpt.yaml'], levels=levels_1sig,
           label=r"EWPO ($1\sigma$)", interpolation_factor=9,
           color='C1')
fpl.contour(**plot_data['global'], levels=levels_123sig,
           label=r"global", interpolation_factor=9,
           color='C3')
plt.axhline(c='0.6', linewidth=1)
plt.axvline(c='0.6', linewidth=1)
plt.xlabel(r"$S$")
plt.ylabel(r"$T$")
plt.legend()
plt.show()
```
4. Conclusions

• Models that explain experimental deviations from the SM in certain observables generically predict also effects in other observables. This is e.g. the case for most models that explain the $B$ anomalies. Consequently, to test such models, one has to consider a global likelihood constructed from as many observables as possible.

• This article shows how to use the python package smelli, which implements a global SMEFT likelihood function. It can be used to either test models, or to interpret data model-independently in the WET and the SMEFT. To date, 399 flavor and other precision observables are included in the likelihood.

• The full global likelihood is work in progress. Since smelli is completely open source, you are welcome to join us on https://github.com/smelli/smelli and to participate in the effort to make smelli truly global.
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