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dave@walnuts:~$ grep -i "eft" /usr/share/dict/british-english
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Left
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dave@walnuts:~$ █
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# Basis construction and translation with DEFT

see B. Gripaos & DS, arXiv:1807.07546, code here

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# Introduction

DEFT takes as input a set of fields and their irreps under a set of  $SU(N)$ -like symmetries. *In principle*, it can output:

- ▶ a list of all the contractions of products of these fields which are invariant under the symmetries, to a given order;
- ▶ a list of the redundancies between these operators (Fierz, IBPs, EOMs, &c.);
- ▶ an arbitrary operator basis;
- ▶ a matrix to convert into and between arbitrary operator bases.

This talk:

1. a look under the hood;
2. the extant obstacles to the practical use of DEFT.

## Irreps of $SU(N)$

All irreps of  $SU(N)$  can be encoded in terms of (anti-)symmetric and traceless combinations of upper and lower indices,  $a, b, \dots = 1, \dots, N$ . The former transforms in the defining irrep., the latter in the conjugate.

$$\begin{array}{ccc} \phi^a \rightarrow U_b^a \phi^b; & \psi_a \rightarrow (U^\dagger)_a^b \psi_b; & A_b^a [A_a^a \equiv 0] \rightarrow \dots \\ N & \bar{N} & N^2 - 1 \end{array}$$

Conjugation raises/lowers indices

$$\phi^a \xleftrightarrow{\text{h.c.}} \phi_a^\dagger$$

There are only three invariant tensors

$$\delta_b^a; \quad \epsilon_{ab\dots z}; \quad \epsilon^{ab\dots z}.$$

## Irreps of Lorentz group

The irreps of  $SO(3, 1)$  are the irreps of  $SU(2)_{\text{lor,L}} \times SU(2)_{\text{lor,R}}$ , encoded by combinations of undotted and dotted indices.

$$e_R^{\dot{\alpha}}, B_{\alpha\beta}$$

Conjugation dots/undots indices

$$\psi_{\alpha} \xleftrightarrow{\text{h.c.}} \psi_{\dot{\alpha}}^{\dagger}$$

More 'conventional' representations can always be repackaged in this form, e.g.,

$$D_{\alpha\dot{\alpha}} = \sigma_{\alpha\dot{\alpha}}^{\mu} D_{\mu},$$
$$F_{\alpha\beta} = 2i(\sigma^{\mu\nu})_{\alpha}^{\gamma} \epsilon_{\gamma\beta} F_{\mu\nu}.$$

# 'Monomial operators'

a product of fields; each field may have some covariant derivatives

In DEFT's internal machinery, all operators are expressed as linear combinations of 'monomial operators'.

$$\delta_{\dot{\gamma}}^{\dot{\alpha}} \delta_{\dot{\delta}}^{\dot{\beta}} \epsilon_{ab} \epsilon^{de} \delta_B^A$$

invariants are L-C epsilons and Kron. deltas

$$H^b \bar{Q}_{LeA}^{\dot{\gamma}} d_R^{\dot{\delta}B} \bar{W}_{\dot{\alpha}\dot{\beta}d}^a$$

$$\bar{W}_d^a \delta_a^d \equiv 0$$

$$\bar{W}_{\dot{\alpha}\dot{\beta}} \epsilon^{\dot{\alpha}\dot{\beta}} \equiv 0$$

Irreps of  $SU(N)$ -like symmetries encoded by upper and lower fundamental indices.

# Lagrangian redundancies

- ▶ Fierz relations  $\epsilon^{\dots}\epsilon_{\dots} = \sum \delta^{\dots}\delta^{\dots}\dots\delta^{\dots}$  (also Schouten identities)

$$\epsilon^{\dot{\alpha}\dot{\beta}}\bar{H}_a e_R^\alpha D_{\alpha\dot{\beta}} D_{\beta\dot{\alpha}} L_L^{\beta a} - \epsilon^{\dot{\alpha}\dot{\beta}}\bar{H}_a e_R^\alpha D_{\beta\dot{\beta}} D_{\alpha\dot{\alpha}} L_L^{\beta a} + \epsilon_{\alpha\beta}\epsilon^{\gamma\delta}\epsilon^{\dot{\alpha}\dot{\beta}}\bar{H}_a e_R^\beta D_{\delta\dot{\beta}} D_{\gamma\dot{\alpha}} L_L^{\alpha a} = 0$$

- ▶ IBPs

$$-\epsilon^{\dot{\alpha}\dot{\beta}}D_{\alpha\dot{\beta}}\bar{L}_{Lb}^{\dot{\gamma}}L_L^{\alpha a}\bar{W}_{\dot{\alpha}\dot{\gamma}a}^b - \epsilon^{\dot{\alpha}\dot{\beta}}\bar{L}_{Lb}^{\dot{\gamma}}D_{\alpha\dot{\beta}}L_L^{\alpha a}\bar{W}_{\dot{\alpha}\dot{\gamma}a}^b + \epsilon^{\dot{\alpha}\dot{\beta}}\bar{L}_{Lb}^{\dot{\gamma}}L_L^{\alpha a}D_{\alpha\dot{\alpha}}\bar{W}_{\dot{\beta}\dot{\gamma}a}^b = 0$$

- ▶ Commuting derivatives

$$-\frac{1}{2}ig'\epsilon^{\alpha\beta}B_{\beta\gamma}D_{\alpha\dot{\alpha}}e_R^\gamma e_R^{\dot{\alpha}} + \frac{1}{2}ig'\epsilon^{\dot{\alpha}\dot{\beta}}\bar{B}_{\dot{\beta}\dot{\gamma}}D_{\alpha\dot{\alpha}}e_R^\alpha e_R^{\dot{\gamma}} - \epsilon^{\alpha\beta}\epsilon^{\dot{\alpha}\dot{\beta}}D_{\gamma\dot{\beta}}D_{\beta\dot{\gamma}}D_{\alpha\dot{\alpha}}e_R^\gamma e_R^{\dot{\gamma}} + \epsilon^{\alpha\beta}\epsilon^{\dot{\alpha}\dot{\beta}}D_{\beta\dot{\gamma}}D_{\gamma\dot{\beta}}D_{\alpha\dot{\alpha}}e_R^\gamma e_R^{\dot{\gamma}} = 0$$

# EOM relations<sup>2</sup>

Local field redefinitions don't affect  $S$ -matrix elements

$$\phi(x) \rightarrow \phi(x) + \frac{U[\phi](x)}{\Lambda^2}$$

$$m_\phi^2 S_2 + S_4 + \frac{S_6}{\Lambda^2} \rightarrow m_\phi^2 S_2 + \left( S_4 + \frac{m_\phi^2}{\Lambda^2} \int dx U \frac{\delta S_2}{\delta \phi} \right) + \frac{1}{\Lambda^2} \left( S_6 + \int dx U \frac{\delta S_4}{\delta \phi} \right) + \mathcal{O}\left(\frac{1}{\Lambda^4}\right)$$

We calculate the **marginal EOM**, whose parts plus the **remainder** we find in higher dimensional terms<sup>1</sup>

$$\begin{aligned} & -\frac{1}{2} \epsilon_{\alpha\beta} \epsilon^{\gamma\delta} \epsilon^{\dot{\alpha}\dot{\beta}} \epsilon_{ab} D_{\delta\dot{\beta}} D_{\gamma\dot{\alpha}} H^b Q_L^{\beta a A} \bar{u}_{R A}^\alpha \\ & + y_e^\dagger \epsilon_{\alpha\beta} \epsilon_{\gamma\delta} \epsilon_{ab} \bar{e}_R^\delta L_L^{\gamma b} Q_L^{\beta a A} \bar{u}_{R A}^\alpha + y_u \epsilon_{\alpha\beta} \epsilon_{\dot{\alpha}\dot{\beta}} \bar{Q}_{L a B}^{\dot{\beta}} Q_L^{\beta a A} \bar{u}_{R A}^\alpha u_R^{\dot{\alpha} B} \\ & + y_d^\dagger \epsilon_{\alpha\beta} \epsilon_{\gamma\delta} \epsilon_{ab} Q_L^{\delta b A} Q_L^{\beta a B} \bar{u}_{R B}^\alpha \bar{d}_{R A}^\gamma - 2\lambda \epsilon_{\alpha\beta} \epsilon_{ab} \bar{H}_d H^b H^d Q_L^{\beta a A} \bar{u}_{R A}^\alpha \end{aligned}$$

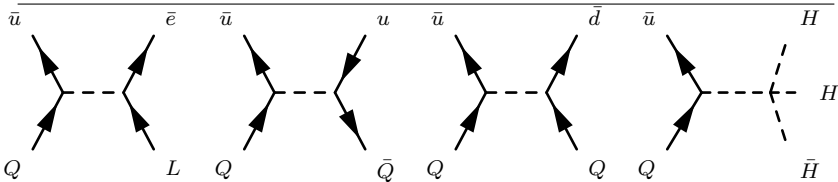
= operators of different dimension, which we ignore

<sup>1</sup>taking care over the **contractions** between the two

<sup>2</sup>Because they look similar, we include Bianchi identities  $D_\mu \tilde{F}^{\mu\nu} = 0$  in this class.

# EOM relations at the amplitude level

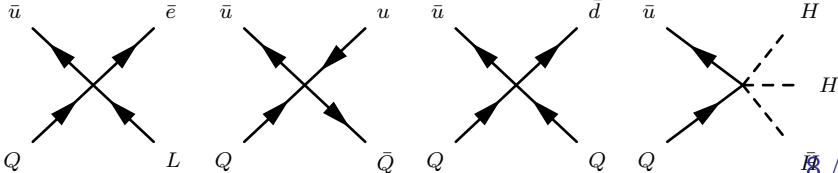
$$Q\bar{u}(-\square - m_H^2)H = Q\bar{u}(y_e^\dagger L\bar{e} + y_u \bar{Q}u + y_d^\dagger Q\bar{d} + 2\lambda H H^\dagger H) + O(\text{dim } 8)$$



with a internal factor of

$$\frac{s}{s - m_H^2} - \frac{m_H^2}{s - m_H^2} = 1$$

is the same as the contact terms





## Put all the relations in a matrix

Each relation is a vector of Wilson coefficients  $\sum_i c_i \mathcal{O}_i = 0$ .  
Collect these as the rows of a matrix.

$$\begin{array}{r} \begin{array}{c} c_1 \\ \downarrow \\ i \\ \leftarrow \text{relation 1} \\ \leftarrow \text{relation 2} \\ \vdots \\ \cdot \\ \cdot \end{array} & \begin{array}{c} c_2 \\ \downarrow \\ 0 \\ 0 \\ \vdots \\ \cdot \\ \cdot \end{array} & \begin{array}{c} c_3 \\ \downarrow \\ y_u \\ 0 \\ \vdots \\ \cdot \\ \cdot \end{array} & \begin{array}{c} \dots \\ 0 \\ 1 \\ \vdots \\ \cdot \\ \cdot \end{array} & \begin{array}{c} c_N \\ \downarrow \\ 0 \\ g_s \\ \vdots \\ \cdot \\ \cdot \end{array} \end{array}$$

The rest is linear algebra, for which we use `sympy`. We've checked the ranks of these matrices for various subsets and simple extensions of the one generation SM, using Hilbert series methods.

## Making a non-redundant basis

1. Perform column operations<sup>3</sup> to express the relations in terms of the coefficients of the operators we want to keep,  $\{\mathcal{B}_1, \mathcal{B}_2\}$ , and the ones we want to eliminate,  $\{\mathcal{D}_1, \mathcal{D}_2, \mathcal{D}_3\}$
2. Perform row operations to reduce to reduced row echelon form

$$\begin{array}{ccccc} c_{\mathcal{D}_1} & c_{\mathcal{D}_2} & c_{\mathcal{D}_3} & c_{\mathcal{B}_1} & c_{\mathcal{B}_2} \\ \downarrow & \downarrow & \downarrow & \downarrow & \downarrow \\ \left( \begin{array}{ccc|cc} 1 & 0 & 0 & * & * \\ 0 & 1 & 0 & * & * \\ 0 & 0 & 1 & * & * \\ \hline 0 & 0 & 0 & 0 & 0 \end{array} \right) \end{array}$$

Iff. the matrix can be put in this form,  $\{\mathcal{B}_1, \mathcal{B}_2\}$  is a valid basis.

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<sup>3</sup>By default, the monomial operators are partially ordered by number of derivatives and epsilon tensors.

# Converting into and between non-redundant bases

The components of the RREF matrix give expressions for any operator  $\mathcal{O}_i \in \{\vec{\mathcal{D}}, \vec{\mathcal{B}}\}$  in terms of the  $\mathcal{B}_i$ :

$$\mathcal{O}_i = R_{ij} \mathcal{B}_j$$

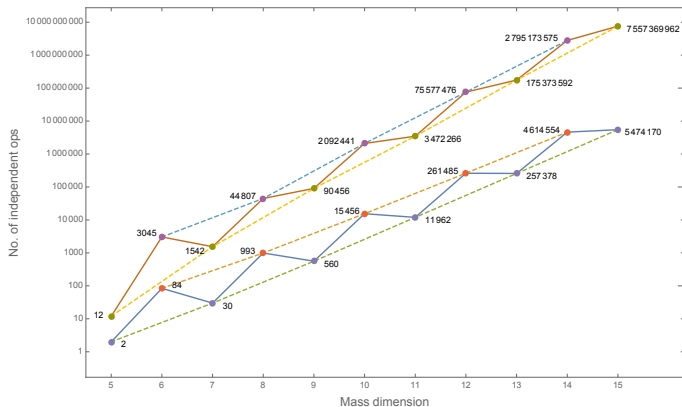
If we can write another basis  $\mathcal{B}'_i$  in terms of the  $\mathcal{O}_i$  (where the  $\mathcal{O}_i$  are by default the monomial terms)

$$\mathcal{B}'_i = S_{ij} \mathcal{O}_j$$

then

$$\mathcal{B}'_i = S_{ij} R_{jk} \mathcal{B}_k$$

# DEFT presently runs out of steam pretty quickly



(From Henning, Lu, Melia, Murayama arXiv:1512.03433)

For the one generation Standard Model on a laptop:

- ▶  $\sim 45$  minutes to generate dim 6 operators and convert from Warsaw to SILH basis
- ▶  $\sim 1$  day to generate a basis of dim 8 operators

## #TODO

- ▶ **Generational indices** Leave them uncontracted or contracted with 0-dimensional spurions in the irreps of the flavour symmetries
- ▶ **Convert between invariants**  $\gamma^\mu, \lambda^A, \dots \leftrightarrow \epsilon, \delta$
- ▶ **Interfaces** Print-out to FeynRules, ... (although note built-in AllYourBases module)
- ▶ **SM in broken phase** Substitute  $B, W, H$  for  $A, Z, W, h$
- ▶ **Transdimensional interactions** Expand EOM relations to higher orders. Systematically calculate corrections to lower order params.
- ▶ **Tabulate higher irreps of  $SU(N)$**  Currently has the rules for (anti)-fundamental, adjoint and all-symmetric tensors
- ▶ **Refactoring** Speed-ups, especially in row reduction methods and memory management

# SM applications?

- ▶ **Matching** Matching a UV theory to the SMEFT is typically done off-shell. One needs to define a ‘Green’s function basis’ and expressions to convert into the usual ‘S-matrix bases’.
- ▶ **Phenomenology of  $d > 6$**  Can study novel processes and define precisely the validity of the EFT expansion.

# Summary

Fields and symmetries in, operators and bases out. In particular, it's been tested against existing results for the one-generation Standard Model.

DEFT is very much a 'proof of concept'. All feedback welcome!