

Using Functional Methods for EFT Calculations

Anders Eller Thomsen

Primarily based on [2412.12270] in collaboration with
J. Fuentes-Martín, A. Moreno-Sánchez, and A. Palavrić

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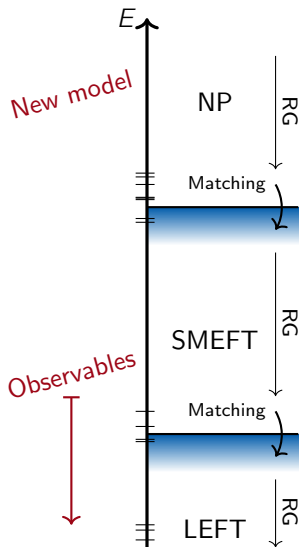
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UNIVERSITÄT
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FOR FUNDAMENTAL PHYSICS

SMEFT Meets ChEFT
TRIUMF, 29 September 2025

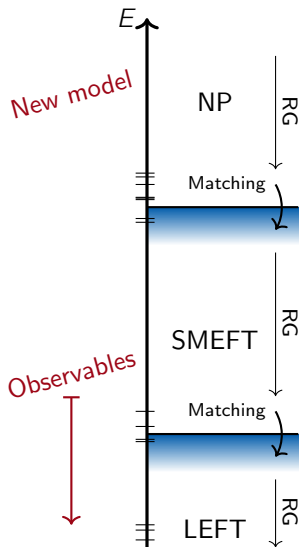


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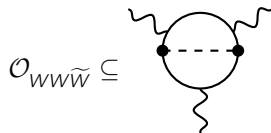
Current status:

- Automated RG flow @ 1-loop
- Automated UV matching @ 1-loop
- Automated SMEFT likelihood
- Calculation of observables @ 1-loop order



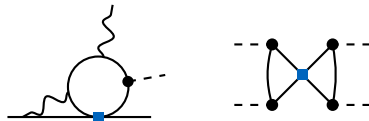
Why we should go beyond one-loop order:

- Restoration of scheme invariance
- New effects at 2-loop order



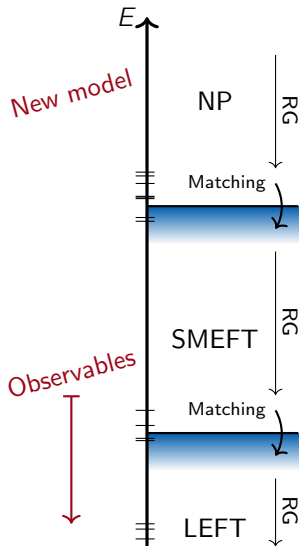
Das Bakshi *et al.* [2103.15861]

- Surprisingly large 2-loop effects



Ardu, Davidson [2103.07212]; Allwicher *et al.* [2302.11584]

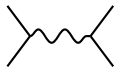
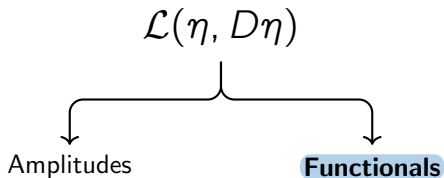
- We might learn something new about QFT



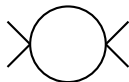
Functional calculations:

- Effective potential (momentum expansion)
- 1-loop **matching calculations**
UOLEA, Matchete,...
- 2-loop RGs with Heat-Kernel
- **2-loop SMEFT RGs**
w.i.p. (see talk by Javier Fuentes-Martín)

What are functional methods anyway?



$$\mathcal{W}[\mathcal{J}]$$



$$\Gamma[\hat{\eta}]$$

Quantum information

Where can we use functional methods?

- EFT matching—extraction of EFT couplings



UV field

$$\sim \int_k \frac{1}{(k^2 - M^2)} \frac{1}{(k+p)^2 - m^2} \sim \frac{-1}{M^2} \int_k \frac{1}{(k+p)^2 - m^2} + \text{local EFT vertex}$$

Changed high- k int

local EFT vertex

Where can we use functional methods?

- EFT matching—extraction of EFT couplings

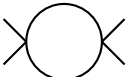


$$\sim \int_k \frac{1}{(k^2 - M^2)} \frac{1}{(k+p)^2 - m^2} \sim \frac{-1}{M^2} \int_k \frac{1}{(k+p)^2 - m^2} + \text{local EFT vertex}$$

UV field

Changed high- k int

- Renormalization/RG functions—extraction of counterterms

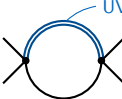


$$\sim \int \frac{d^4 k}{(2\pi)^4} \frac{1}{(k^2 - M^2)^2} \longrightarrow \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k^2 - M^2)^2} + \delta\lambda$$

local UV counterterm

Where can we use functional methods?

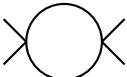
- EFT matching—extraction of EFT couplings



$$\sim \int_k \frac{1}{(k^2 - M^2)} \frac{1}{(k+p)^2 - m^2} \xrightarrow{\text{Changed high-}k \text{ int}} \frac{-1}{M^2} \int_k \frac{1}{(k+p)^2 - m^2} + \text{local EFT vertex}$$

The diagram shows a bubble with two external lines. A blue arc above the bubble is labeled "UV field". The integral is shown with a blue box around the $\frac{1}{(k^2 - M^2)}$ term. A blue arrow points from this term to the $\frac{-1}{M^2}$ term, labeled "Changed high- k int". The final term is a blue circle with four external lines, labeled "local EFT vertex".

- Renormalization/RG functions—extraction of counterterms



$$\sim \int \frac{d^4 k}{(2\pi)^4} \frac{1}{(k^2 - M^2)^2} \longrightarrow \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k^2 - M^2)^2} + \delta\lambda$$

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- Amplitudes? Why not?!

But there might be little point... and some *trouble* $\leftrightarrow$ non-local

Where can we use functional methods?

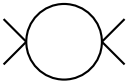
- EFT matching—extraction of EFT couplings



$$\sim \int_k \frac{1}{(k^2 - M^2)} \frac{1}{(k+p)^2 - m^2} \xrightarrow{\text{Changed high-}k \text{ int}} \frac{-1}{M^2} \int_k \frac{1}{(k+p)^2 - m^2} + \text{local EFT vertex}$$

The diagram shows a bubble with two external lines. A blue arrow points to the top arc of the bubble, labeled "UV field". The integral is shown with a blue box around the $\frac{1}{(k^2 - M^2)}$ term. A blue arrow points from this term to the $\frac{-1}{M^2}$ term, labeled "Changed high- k int". The final term is a blue circle with four external lines, labeled "local EFT vertex".

- Renormalization/RG functions—extraction of counterterms



$$\sim \int \frac{d^4 k}{(2\pi)^4} \frac{1}{(k^2 - M^2)^2} \longrightarrow \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k^2 - M^2)^2} + \delta\lambda$$

The diagram shows a bubble with two external lines. The integral is shown with a blue box around the $\frac{1}{(k^2 - M^2)^2}$ term. The final term is a blue $\delta\lambda$, labeled "local UV counterterm".

- Amplitudes? Why not?!

But there might be little point... and some *trouble* $\leftrightarrow$ non-local

But what exactly are the advantages of functional methods?

A direct approach to RG quantities

Diagrammatic approach (tensor quantity):

$$\begin{aligned} \text{Diagram 1: Circle with '1PI' and four external lines} &= \int dx_1 e^{-ip_1 \cdot x_1} \dots \frac{\delta}{\delta \hat{\eta}(x_1)} \dots \Gamma[\hat{\eta}] \Big|_{\hat{\eta}=0} \\ \text{Diagram 2: Blue dot with four external lines} &= \int dx_1 e^{-ip_1 \cdot x_1} \dots \frac{\delta}{\delta \hat{\eta}(x_1)} \dots \delta S[\hat{\eta}] \Big|_{\hat{\eta}=0} \end{aligned}$$

A direct approach to RG quantities

Diagrammatic approach (tensor quantity):

$$\begin{aligned} \text{1PI} &= \int dx_1 e^{-ip_1 \cdot x_1} \dots \frac{\delta}{\delta \hat{\eta}(x_1)} \dots \Gamma[\hat{\eta}] \Big|_{\hat{\eta}=0} \\ \text{Cross} &= \int dx_1 e^{-ip_1 \cdot x_1} \dots \frac{\delta}{\delta \hat{\eta}(x_1)} \dots \delta S[\hat{\eta}] \Big|_{\hat{\eta}=0} \end{aligned}$$

Functional method (scalar quantity):

- EFT matching

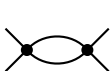
$$S_{\text{EFT}}[\hat{\phi}] \subseteq \Gamma_{\text{UV}}[\hat{\eta}] \quad \longrightarrow \quad S_{\text{EFT}} = \overset{\text{expansion operator}}{\mathcal{T}}_{\text{hard}} \Gamma_{\text{UV}}$$

- Renormalization

$$-\delta S[\hat{\eta}] \subseteq \Gamma[\hat{\eta}] \quad \longrightarrow \quad \delta S = -\mathcal{T}_\delta \Gamma$$

No repetition from symmetries

Diagrammatic approach: multiple permutations of external lines:

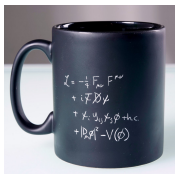


Functional method: contributions exclusively determined by internal lines

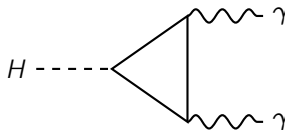


Saving more pronounced with more external lines or less internal symmetry, e.g., W^3 , $(H^* H^*)^3$, $\phi^6, \dots$

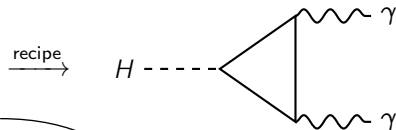
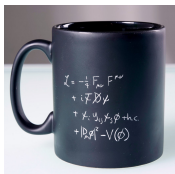
Benefits from a master formula



recipe
→



Benefits from a master formula



$$\sum_{\text{1-loop}} \text{diagram} = \frac{i\hbar}{2} \text{STr} \log Q \xrightarrow{\text{recipe}} \mathcal{L}_{\text{LEFT}}$$

- **Bookkeeping:** organization by topology/internal propagators
- **Manifest covariance:** There is no need to separate out

$$D_\mu = \partial_\mu - iA_\mu$$

- **Recycling:** the methods recycle known loop-integral techniques

Functional calculation step-by-step

General formulas:

1. Expand quantum effective action (background field method)
2. Cast topologies as covariant momentum integrals
3. Determine expansion operator T_{uv} (what part of Γ to evaluate)

Computer implementation:

4. Expand the super diagram:
 - Expand kinetic operators by spin (scalar, fermion, vector)
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5. Perform tadpole momentum integrals with known techniques
6. Simplify resultant Lagrangian

Semi-classical approximation

The **quantum effective action** is the generating functional of 1PI functions:

$$e^{i\Gamma_{\text{UV}}[\hat{\eta}]} = \int \mathcal{D}\eta \exp\left(iS_{\text{UV}}[\eta + \hat{\eta}] + iJ_I \eta^I\right), \quad \frac{\delta\Gamma[\hat{\eta}]}{\delta\hat{\eta}^I} = -J_I$$

$\searrow \int_x J_a(x) \eta^a(x)$

The action is expanded by legs and loop order (including counterterms):

$$S[\eta + \hat{\eta}] = \hat{S} + \sum_{\ell=0}^{\infty} \hbar^{\ell} \left(\eta_I \hat{\mathcal{V}}_I^{(\ell)} + \frac{1}{2} \eta_I \hat{\mathcal{V}}_{IJ}^{(\ell)} \eta_J + \frac{1}{6} \eta_I \eta_J \eta_K \hat{\mathcal{V}}_{IJK}^{(\ell)} + \frac{1}{24} \eta_I \eta_J \eta_K \eta_L \hat{\mathcal{V}}_{IJKL}^{(\ell)} + \dots \right)$$

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With a **saddlepoint approximation**

$$e^{i\Gamma_{\text{UV}}[\hat{\eta}]} = e^{i\hat{S}} \int \mathcal{D}\eta \exp\left(\frac{i}{2} \eta^I \hat{\mathcal{Q}}_{IJ} \eta^J\right) \left[1 + \frac{i}{24} \eta_I \eta_J \eta_K \eta_L \hat{\mathcal{V}}_{IJKL}^{(0)} - \frac{1}{72} (\eta_I \eta_J \eta_K \hat{\mathcal{V}}_{IJK}^{(0)})^2 + \dots \right]$$

$\nearrow \mathcal{V}_{IJ}^{(0)}$

Semi-classical approximation

The loop-expansion is a polynomial of tensor contractions

$$\begin{aligned} \Gamma[\hat{\eta}] = & \hat{S} + \frac{i\hbar}{2} \overbrace{\zeta_{II} (\log \hat{Q})_{II}}^{\text{STr log } Q} + \frac{i\hbar^2}{2} \zeta_{II} \hat{Q}_{IJ}^{-1} \hat{V}_{JI}^{(1)} - \frac{\hbar^2}{8} \hat{Q}_{IJ}^{-1} \hat{Q}_{KL}^{-1} \hat{V}_{IJKL}^{(0)} \\ & + \frac{\hbar^2}{12} \zeta_{IM} \zeta_{JN} \zeta_{IN} \hat{Q}_{LI}^{-1} \hat{Q}_{MJ}^{-1} \hat{Q}_{NK}^{-1} \hat{V}_{IJK}^{(0)} \hat{V}_{LMN}^{(0)} + \mathcal{O}(\hbar^3) \end{aligned}$$

Spin statistics: ± 1

or as supergraphs

$$\Gamma_{uv} = S_{uv} + \frac{i}{2} \log \text{[circle]} + \frac{i}{2} \text{[circle with top vertex]}^{(1)} - \frac{1}{8} \text{[two circles connected at bottom]} + \frac{1}{12} \text{[circle with two vertices]} + \mathcal{O}(\hbar^3)$$

Semi-classical approximation

The loop-expansion is a polynomial of tensor contractions

$$\Gamma[\hat{\eta}] = \hat{S} + \frac{i\hbar}{2} \overbrace{\zeta_{II} (\log \hat{Q})_{II}}^{\text{STr log } Q} + \frac{i\hbar^2}{2} \zeta_{II} \hat{Q}_{IJ}^{-1} \hat{V}_{JI}^{(1)} - \frac{\hbar^2}{8} \hat{Q}_{IJ}^{-1} \hat{Q}_{KL}^{-1} \hat{V}_{IJKL}^{(0)} \\ + \frac{\hbar^2}{12} \zeta_{IM} \zeta_{JN} \zeta_{IN} \hat{Q}_{LI}^{-1} \hat{Q}_{MJ}^{-1} \hat{Q}_{NK}^{-1} \hat{V}_{IJK}^{(0)} \hat{V}_{LMN}^{(0)} + \mathcal{O}(\hbar^3)$$

Spin statistics: ± 1

or as supergraphs

$$\Gamma_{uv} = S_{uv} + \frac{i}{2} \log \text{ (loop with one vertex) } + \frac{i}{2} \text{ (loop with one vertex and one external line) } - \frac{1}{8} \text{ (two loops) } + \frac{1}{12} \text{ (triangle) } + \mathcal{O}(\hbar^3)$$

All **vertices and propagators are fully dressed** with background fields

$$Q_{IJ}[\hat{\eta}] = \frac{\delta^2 S^{(0)}}{\delta \eta^I \delta \eta^J}[\hat{\eta}] = \left(\text{diagram of a loop with two external lines} \right)^{-1}$$

$$\mathcal{V}_{IJK}^{(0)}[\hat{\eta}] = \text{diagram of a vertex with three external lines}$$

The functional index

$$I = x, \left\{ \begin{array}{ll} \text{scalar} : () & \left\{ \begin{array}{l} H : (i) \\ H^* : (i) \end{array} \right. \\ \text{vector} : (\mu) & \left\{ \begin{array}{l} B : () \\ \vdots \end{array} \right. \\ \text{spinor} : (\alpha) & \left\{ \begin{array}{l} q : (c, i, p) \\ q^C : (c, i, p) \\ \vdots \end{array} \right. \\ \text{(anti-)ghost} : () & \left\{ \begin{array}{l} \vdots \end{array} \right. \end{array} \right.$$

The functional index

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Individual diagrams are embedded in supergraphs

$$\hat{\mathcal{Q}}_{LI}^{-1} \hat{\mathcal{Q}}_{MJ}^{-1} \hat{\mathcal{Q}}_{NK}^{-1} \hat{\mathcal{V}}_{IJK}^{(0)} \hat{\mathcal{V}}_{LMN}^{(0)} = \text{diagram 1} \supseteq \text{diagram 2} \supseteq \text{diagram 3}$$

1:1 correspondence fails for external gauge bosons

Functional calculation step-by-step

General formulas:

1. Expand quantum effective action (background field method)
2. Cast topologies as covariant momentum integrals
3. Determine expansion operator $\mathcal{T}_{UV}$ (what part of Γ to evaluate)

Computer implementation:

4. Expand the super diagram:
 - Expand kinetic operators by spin (scalar, fermion, vector)
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Covariant functional Feynman rules

Locality of the action: functional tensors factors as differential operators

$$\mathcal{V}_{l_1 \dots l_n} = \sum_{m_2 \dots m_n} V_{a_1 a'_2 \dots a'_n}^{(m_2, \dots, m_n)}(x_1) \prod_{i=2}^n P_{x_i}^{m_i} \delta_{a'_i a_i}(x_1, x_i)$$

$$Q_{IJ} = \underbrace{Q_{ac}(x, P_x)}_{\text{manifestly covariant}} \delta^c_b(x, y)$$

iD_μ

Covariant functional Feynman rules

Locality of the action: functional tensors factors as differential operators

$$\mathcal{V}_{l_1 \dots l_n} = \sum_{m_2 \dots m_n} V_{a_1 a'_2 \dots a'_n}^{(\underline{m}_2, \dots, \underline{m}_n)}(x_1) \prod_{i=2}^n P_{x_i}^{\underline{m}_i} \delta_{a'_i a_i}(x_1, x_i)$$
$$\mathcal{Q}_{IJ} = Q_{ac}(x, P_x) \overset{iD_\mu}{\delta^c_b}(x, y)$$

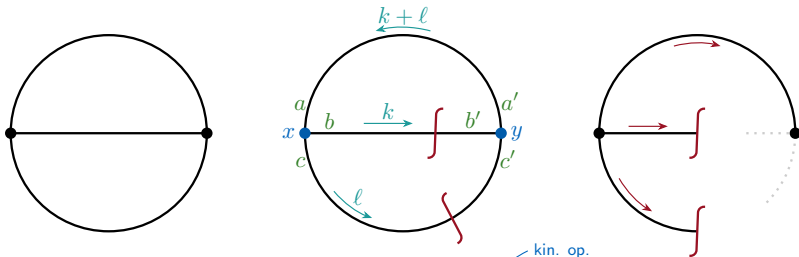
And covariance is manifest with the **covariant delta functions**

$$\frac{\delta \eta^a(x)}{\delta \eta^b(y)} = \delta^a_b(x, y), \quad \delta^a_b(x, y) \equiv \overset{\text{Wilson line}}{U^a_b}(x, y) \delta(x - y)$$
$$D_x^\mu \delta^a_b(x, y) = -D_y^\mu \delta^a_b(x, y)$$

The **background gauge field is a connection** in the field manifold

$$f^a(x) = U^a_b(x, y) f^b(y), \quad U(x, y) = \mathcal{P} \exp \left[i \int_x^y dz^\mu \hat{A}_\mu(z) \right]$$

Sunset formula



$$E_{a_i b_j}^{(\underline{m}_i, \underline{n}_j)}(x, P_x + k) \equiv (-1)^{m_i} (P_x + k)^{\underline{m}_i} Q_{a_i b_j}^{-1}(x, P_x + k) (P_x + k)^{\underline{n}_j}$$

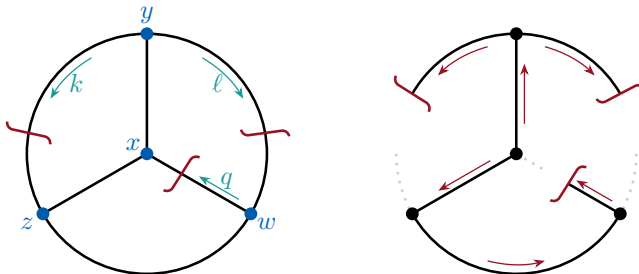
Sunset master formula:

$$\frac{\hbar^2}{12} \sum_{m_i, n_i} \iint_x \iint_{kl} E_{aa'}^{(0,0)}(y, P_y - k - \ell) V_{abc}^{(\underline{m}_1, \underline{n}_1)}(x) V_{a'b'c'}^{(\underline{m}_2, \underline{n}_2)}(y) \\ \times [E^{(\underline{m}_1, \underline{m}_2)}(x, P_x + k) U(x, y)]_{bb'} [E^{(\underline{n}_1, \underline{n}_2)}(x, P_x + \ell) U(x, y)]_{cc'} \Big|_{y=x}$$

Grassmanian factor: $\zeta_{aa'} \zeta_{bb'} \zeta_{cc'} \zeta_{ab} \zeta_{ac} \zeta_{a'b} \zeta_{a'c} \zeta_{bc'} \zeta_{a'b'} \zeta_{a'c'} = \pm 1$

Tetrahedron formula

Systematic approach to deriving master formulas from diagram topologies:



$$\begin{aligned}
 & \frac{i\hbar^3}{24} \sum_{m_i, n_i} \int_x \int_{k\ell q} E_{a_1 b_1}^{(0,0)}(y, P_y + k + \ell) E_{a_2 c_1}^{(\underline{m}_1, 0)}(z, P_z + q - k - \ell) E_{c_2 d_3}^{(\underline{m}_3, \underline{n}_4)}(w, P_w + q - \ell) \\
 & \times V_{a_1 a_2 a_3}^{(\underline{m}_1, \underline{n}_1)}(x) V_{b_1 b_2 b_3}^{(\underline{m}_2, \underline{n}_2)}(y) V_{c_1 c_2 c_3}^{(\underline{m}_3, \underline{n}_3)}(z) V_{d_1 d_2 d_3}^{(\underline{m}_4, \underline{n}_4)}(w) [E^{(\underline{m}_2, \underline{n}_3)}(y, P_y + k) U(y, z)]_{b_2 c_3} \\
 & \times [E^{(\underline{n}_2, \underline{m}_4)}(y, P_y + \ell) U(y, w)]_{b_3 d_2} [E^{(0, \underline{n}_1)}(w, P_w + q) U(w, x)]_{d_1 a_3} \Big|_{y, z, w=x}
 \end{aligned}$$

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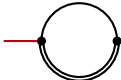
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Expansion by regions

With expansion by regions we can **separate scales in loop integrals**, e.g., a 2-point function with $p^2, m^2 \ll \Lambda^2$:

$$I = \text{diagram} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k+p)^2 - m^2} \frac{1}{k^2 - \Lambda^2}$$
A Feynman diagram representing a 2-point function. It consists of a circular loop with two external red lines attached to the left and right vertices. The loop is drawn with two concentric circles, indicating a specific regularization scheme.

Expansion by regions

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$$I = \text{[bubble diagram]} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k+p)^2 - m^2} \frac{1}{k^2 - \Lambda^2}$$
$$I_h = \text{[hard region diagram]} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{k^2} \frac{1}{k^2 - \Lambda^2} + \dots$$

$k^2 \gtrsim \Lambda^2$

Expansion by regions

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$$\begin{aligned} I &= \text{[bubble diagram]} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k+p)^2 - m^2} \frac{1}{k^2 - \Lambda^2} \\ I_h &= \text{[hard diagram]} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{k^2} \frac{1}{k^2 - \Lambda^2} + \dots \\ I_s &= \text{[soft diagram]} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k+p)^2 - m^2} \frac{-1}{\Lambda^2} + \dots \end{aligned}$$

$k^2 \ll \Lambda^2$

Expansion by regions

With expansion by regions we can **separate scales in loop integrals**, e.g., a 2-point function with $p^2, m^2 \ll \Lambda^2$:

$$\begin{aligned} I &= \text{diagram of a bubble loop} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k+p)^2 - m^2} \frac{1}{k^2 - \Lambda^2} \\ I_h &= \text{diagram of a hard region (blue circle)} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{k^2} \frac{1}{k^2 - \Lambda^2} + \dots \\ I_s &= \text{diagram of a soft region (red circle)} = \int \frac{d^d k}{(2\pi)^d} \frac{1}{(k+p)^2 - m^2} \frac{-1}{\Lambda^2} + \dots \end{aligned}$$

In dim. reg. integrals equal the sum of their **hard** and **soft** regions

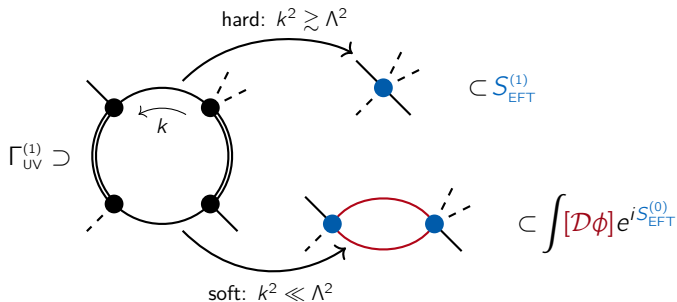
Beneke, Smirnov [hep-ph/9711391]; Jantzen [1111.2589]

$$I = I_h + I_s$$

The regions I_h and I_s are **systematically improvable** power series in $1/\Lambda^2$

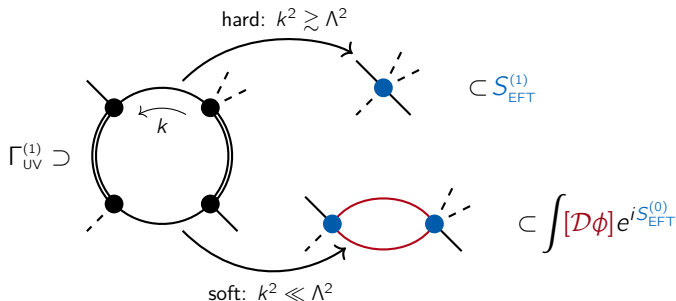
Separation of scales in one-loop matching

Mixed (heavy–light) loop example:



Separation of scales in one-loop matching

Mixed (heavy–light) loop example:



- $\Gamma_{UV}^{(1)}|_{\text{soft}}$: long-distance contributions included in 1-loop matrix elements of tree-level EFT operators

$$\Gamma_{UV}^{(1)}|_{\text{soft}} = \Gamma_{EFT}^{(1)}$$

- $\Gamma_{UV}^{(1)}|_{\text{hard}}$: short-distance contributions going into the EFT operators

Fuentes-Martin *et al.* [1607.02142]; Zhang [1610.00710]

$$\Gamma_{UV}^{(1)}|_{\text{hard}} = S_{EFT}^{(1)}$$

Hard-region matching formula

$$S_{\text{EFT}}[\phi] = \Gamma_{\text{UV}}[\hat{\Phi}, \phi]_{\text{hard}}, \quad \frac{\delta \Gamma_{\text{UV}}|_{\text{hard}}}{\delta \Phi}[\hat{\Phi}, \phi] = 0$$

“hard” denotes the part without *any* soft loop momenta (it includes all tree-level contributions)

Fuentes-Martín, Moreno-Sánchez, Palavrić, *AET* [2311.13630] [2412.12270]

See also Fuentes-Martín *et al.* [1607.02142]; Zhang [1610.00710]

- Allows for matching without calculating any EFT loops
- The expansion operator $T_{\text{UV}} = T_{\text{hard}}$ selects a local part of Γ_{UV} to evaluate
- Proven to all orders in the loop expansion

Functional calculation step-by-step

General formulas:

1. Expand quantum effective action (background field method)
2. Cast topologies as covariant momentum integrals
3. Determine expansion operator T_{UV} (what part of Γ to evaluate)
 - For matching calculations, $T_{UV} = T_{\text{hard}}$
 - For RGE calculations, $T_{UV} = \bar{R}^*$

Computer implementation:

4. Expand the super diagram:
 - Expand kinetic operators by spin (scalar, fermion, vector)
 - Expand propagators in the covariant momentum operators
 - Sum over possible field labels $\rightsquigarrow$ substitute field-specific interactions
5. Perform tadpole momentum integrals with known techniques
6. Simplify resultant Lagrangian

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Evaluating momentum-space topologies

E.g., the sunset formula:

$$\frac{\hbar^2}{12} \sum_{m_i, n_i} \int_x \int_{k\ell} E_{aa'}^{(0,0)}(y, P_y - k - \ell) V_{abc}^{(\underline{m}_1, \underline{n}_1)}(x) V_{a'b'c'}^{(\underline{m}_2, \underline{n}_2)}(y) \\ \times [E^{(\underline{m}_1, \underline{m}_2)}(x, P_x + k) U(x, y)]_{bb'} [E^{(\underline{n}_1, \underline{n}_2)}(x, P_x + \ell) U(x, y)]_{cc'} \Big|_{y=x}$$

RG calculations with T_{UV} (k, ℓ large) call for expansion of the propagators:

$$T_{UV} E^{(\underline{m}, \underline{n})} \supset T_{UV} Q^{-1} = T_{UV} \Delta^{-1} \sum_{n=0}^{\infty} (X \Delta^{-1})^n$$

- Typical propagator: $\Delta^{-1} \sim \frac{1}{(k + P_x)^2 - m^2} = \frac{1}{k^2} - \frac{2k \cdot P_x}{k^4} + \dots$
- Typical X -term: $X_{\phi A_\mu} \supset 2g\phi P_x^\mu$
- Closed formula: $P_x^{\mu_1} \dots P_x^{\mu_m} U(x, y) \Big|_{y=x} = p_m(P_x, G_{\mu\nu})$ polynomial

Functional calculation step-by-step

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6. **Simplify resultant Lagrangian** $\rightsquigarrow$ challenging in generality (harder than “just” determining an EFT basis).

Summary: **tensor contract; expand; loop-integrate**

Functional calculation step-by-step

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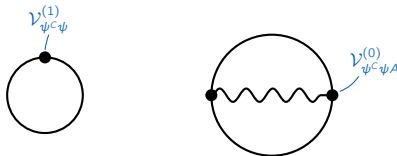
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Example: Euler–Heisenberg

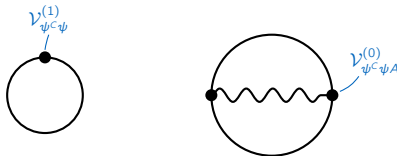
Let's integrate out the electron at 2-loop orders!



2-loop, EFT-order 8, by-hand calculation

Example: Euler–Heisenberg

Let's integrate out the electron at 2-loop orders!



2-loop, EFT-order 8, by-hand calculation

$$\begin{aligned}
 \mathcal{L}_{\text{EFT}} = & -\frac{1}{4e^2} \left[1 + \frac{e^2}{16\pi^2} \frac{4}{3} \log \frac{\mu^2}{m^2} + \frac{e^4}{(16\pi^2)^2} \left(\frac{13}{3} - 4 \log \frac{\mu^2}{m^2} \right) \right] F^{\mu\nu} F_{\mu\nu} \\
 & + \left[\frac{1}{16\pi^2} \frac{7}{90} \frac{1}{m^4} - \frac{1}{(16\pi^2)^2} \frac{e^2}{m^4} \left(\frac{14}{15} \log \frac{\mu^2}{m^2} + \frac{247}{810} \right) \right] (F^{\mu\nu} F_{\nu\sigma})^2 \\
 & - \left[\frac{1}{16\pi^2} \frac{1}{36} \frac{1}{m^4} - \frac{1}{(16\pi^2)^2} \frac{e^2}{m^4} \left(\frac{1}{3} \log \frac{\mu^2}{m^2} - \frac{49}{324} \right) \right] (F^{\mu\nu} F_{\mu\nu})^2
 \end{aligned}$$

agrees with Grozin [1212.5144]

consistent with μ -independence

Fuentes-Martin, Moreno-Sánchez, Palavrić, AET [2412.12270]

Summary

- Functional methods are an excellent match for **all RG calculations**
- The parallels with ordinary Feynman diagrams ensure that we may borrow known integrals techniques
- The techniques are extendable to 3-loop orders and beyond (subject to the usual computational limitations)
- Other interesting results: all-order EFT matching formula
- Practical implementation of the techniques [see talk by Javier Fuentes-Martín](#)

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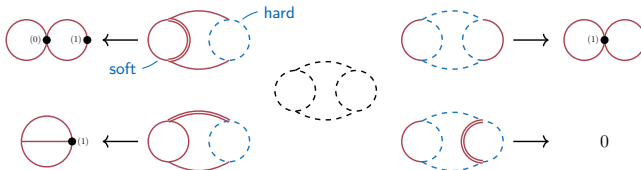
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Thank you!

Backup

Proof of hard-region matching formula

- Expansion by regions (soft/hard) extended to arbitrary loop order
- Compare UV and EFT effective action (**induction in loop order**)
 - Decompose UV topologies (figure is not comprehensive)



- Use EFT vertices produced from lower order matching

$$i\widehat{\mathcal{W}}_{ijk}^{(\ell)} = i \text{---} \text{blue circle} \begin{matrix} \nearrow j \\ \searrow k \end{matrix} + \sum_{\text{perm.}} i \text{---} \text{blue circle} \text{---} \text{red dot} \begin{matrix} \nearrow j \\ \searrow k \end{matrix} + \sum_{\text{perm.}} \text{blue circle} \text{---} \text{red line} \text{---} \text{red dot} \begin{matrix} \nearrow j \\ \searrow k \end{matrix} + \text{blue circle} \text{---} \text{red line} \text{---} \text{red dot} \begin{matrix} \nearrow i \\ \searrow j \\ \searrow k \end{matrix}$$

- Account for combinatorics and symmetry factors

See Fuentes-Martín, Moreno-Sánchez, Palavrić, **AET** [2412.12270]