

Package-X 2.1 and CollierLink

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ACAT Track 3

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INTERACTIONS

What is *Package-X* ?

Released in 2015 :

Package-X is a *Mathematica* package to (mainly) calculate one-loop integrals analytically.

Calculations carried out in 3 steps:

$$\int \frac{d^d k}{(2\pi)^d} \frac{k^\mu k^\nu}{[(k+p)^2 - m^2] k^2} \text{ at } p^2 = 0$$

1. `LoopIntegrate[kμ kν, k, {k+p, m}, {k, 0}]`
`pμ pν PVB[0, 2, p.p, 0, m] + gμ,ν PVB[1, 0, p.p, 0, m]`

2. `% /. p.p → 0`
`pμ pν PVB[0, 2, 0, 0, m] + gμ,ν PVB[1, 0, 0, 0, m]`

3. `% // LoopRefine`
$$\left(\frac{1}{9} + \frac{1}{3} \left(\frac{1}{\epsilon} + \text{Log} \left[\frac{\mu^2}{m^2} \right] \right) \right) p_\mu p_\nu + \left(\frac{3m^2}{8} + \frac{1}{4} m^2 \left(\frac{1}{\epsilon} + \text{Log} \left[\frac{\mu^2}{m^2} \right] \right) \right) g_{\mu,\nu}$$

Important goal: ***EASY TO USE***

What is *Package-X* ?

Version 2 Released in 2016 :

A) boxes

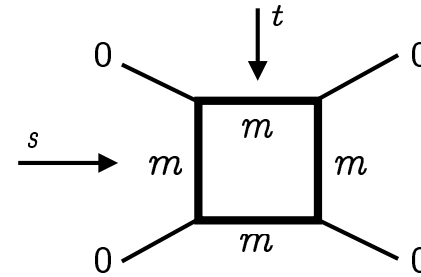
```
LoopIntegrate[1, k, {k, m}, {k + p1, m},
  {k - p3 + p1, m}, {k - p2, m}]
```

```
PVD[0, 0, 0, 0, p1.p1, p3.p3,
  p1.p1 + 2 p1.p2 - 2 p1.p3 + p2.p2 - 2 p2.p3 + p3.p3,
  p2.p2, p1.p1 - 2 p1.p3 + p3.p3,
  p1.p1 + 2 p1.p2 + p2.p2, m, m, m, m]
```

```
LoopRefine[%] // D0Expand
```

$$\text{ConditionalExpression}\left[\frac{1}{s \sqrt{1 - \frac{4m^2}{s} - \frac{4m^2}{t}}} 2 \left(-\frac{\pi^2}{2} - \text{Log}\left[\frac{\left(1 + \sqrt{1 - \frac{4m^2}{s}}\right) t \left(\sqrt{1 - \frac{4m^2}{s}} - \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right]^2 - \text{Log}\left[\frac{s \left(1 + \sqrt{1 - \frac{4m^2}{t}}\right) \left(\sqrt{1 - \frac{4m^2}{t}} - \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right]^2 + \right. \\ \left. \text{Log}\left[-\frac{t \left(\sqrt{1 - \frac{4m^2}{s}} - \sqrt{1 - \frac{4m^2}{st}}\right)^2}{4m^2}\right] \text{Log}\left[-\frac{s \left(\sqrt{1 - \frac{4m^2}{t}} - \sqrt{1 - \frac{4m^2}{st}}\right)^2}{4m^2}\right] + 2 \text{Log}\left[-\frac{s \left(\sqrt{1 - \frac{4m^2}{s}} + \sqrt{1 - \frac{4m^2}{st}}\right) \left(-\sqrt{1 - \frac{4m^2}{t}} + \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right]^2 - \right. \\ \left. 2 \text{PolyLog}\left[2, -\frac{\left(-1 + \sqrt{1 - \frac{4m^2}{s}}\right) s \left(\sqrt{1 - \frac{4m^2}{s}} - \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right] + 2 \text{PolyLog}\left[2, \frac{\left(-1 + \sqrt{1 - \frac{4m^2}{s}}\right) t \left(\sqrt{1 - \frac{4m^2}{s}} - \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right] + \right. \\ \left. 2 \text{PolyLog}\left[2, \frac{s \left(-1 + \sqrt{1 - \frac{4m^2}{t}}\right) \left(\sqrt{1 - \frac{4m^2}{t}} - \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right] - 2 \text{PolyLog}\left[2, -\frac{\left(-1 + \sqrt{1 - \frac{4m^2}{t}}\right) t \left(\sqrt{1 - \frac{4m^2}{t}} - \sqrt{1 - \frac{4m^2}{st}}\right)}{4m^2}\right] \right], st (st - 4m^2 (s+t)) > 0]$$

A.I.Davydychev
arXiv: 9307323



What is *Package-X* ?

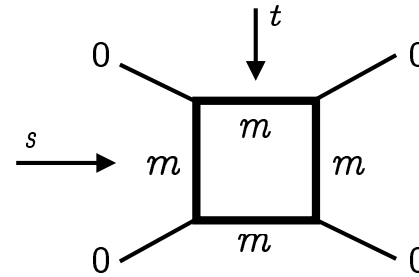
Version 2 Released in 2016 :

B) series expansions

Large mass expansion ($m \sim \infty$)

`LoopRefineSeries[%%, {m, Infinity, 12}]`

$$\frac{1}{6 m^4} + \frac{s+t}{60 m^6} + \frac{2 s^2 + s t + 2 t^2}{840 m^8} + \frac{3 s^3 + s^2 t + s t^2 + 3 t^3}{7560 m^{10}} + \frac{12 s^4 + 3 s^3 t + 2 s^2 t^2 + 3 s t^3 + 12 t^4}{166320 m^{12}} + O\left[\frac{1}{m}\right]^{13}$$



Forward limit ($t \sim 0$)

`LoopRefineSeries[%%, {t, 0, 2}]`

$$\left(\frac{2}{m^2 s} + \frac{\text{DiscB}[s, m, m]}{m^2 s} \right) + \left(-\frac{4(3 m^2 - s)}{9 m^4 s^2} - \frac{(4 m^2 - s) \text{DiscB}[s, m, m]}{6 m^4 s^2} \right) t + \left(\frac{240 m^4 - 140 m^2 s + 23 s^2}{225 m^6 s^3} + \frac{(4 m^2 - s)^2 \text{DiscB}[s, m, m]}{30 m^6 s^3} \right) t^2 + O[t]^3$$

C) improved numerics

machine precision `ScalarD0[5.4, 1.2, 5.4, 1.2, 54.1, -12, 1.1, 1.1, 1.1, 2.3]`

precision `-0.0174954 + 0.0101058 i`

arbitrary precision `ScalarD0[5.4`20, 1.2`20, 5.4`20, 1.2`20, 54.1`20, -12`20, 1.1`20, 1.1`20, 1.1`20, 2.3`20]`

precision `-0.0174953803351620551279058 + 0.01010581673617592175607057 i`

What is *Package-X* ?

Version 2 Released in 2016 :

D) Dirac algebra

Calculate traces (already in v1.0)

```
Spur[k.γ + m 1, γμ, (k + p).γ + m 1, γν]
```

```
8 kμ kν + 4 kν pμ + 4 kμ pν + 4 m2 gμν - 4 k.k gμν - 4 k.p gμν
```

Put open fermion lines in canonical form:

```
⟨u[p2, m], p1.γ, γμ, p2.γ, u[p1, m]⟩ // FermionLineExpand
```

```
-2 i m ⟨u[p2, m], σμ,{-p1+p2}, u[p1, m]⟩ + ⟨u[p2, m], γμ, u[p1, m]⟩ (3 m2 - 2 p1.p2)
```

Compute a Feynman diagram (QED vertex function):

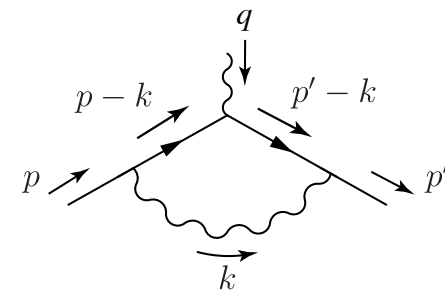
```
LoopIntegrate[⟨u[p', m], γν, (p'.γ - k.γ + m 1), γμ,
  (p.γ - k.γ + m 1), γν, u[p, m]⟩, k, {k - p', m}, {k - p, m},
  {k, 0, 1}] /. {p.p → m2, p'.p' → m2, p.p' → -q.q / 2 + m2} //
```

```
LoopRefine
```

```
2 i m DiscB[q.q, m, m] ⟨u[p', m], σμ,{-p+p'}, u[p, m]⟩
-----
4 m2 - q.q
```

```
⟨u[p', m], γμ, u[p, m]⟩ ⎛ -  $\frac{\text{DiscB}[q.q, m, m] (8 m^2 - 3 q.q)}{4 m^2 - q.q}$  +
```

```
⎛ 1 -  $\frac{2 \text{DiscB}[q.q, m, m] (2 m^2 - q.q)}{4 m^2 - q.q}$  ⎞ ⎛  $\frac{1}{\epsilon} + \text{Log}\left[\frac{\mu^2}{m^2}\right]$  ⎞ + 2 (2 m2 - q.q) ScalarC0IR6[q.q, m, m]
```



What is *Package-X* ?

Version 2 Released in 2016 :

E) Intuitive and readable user interface

Input and Output in StandardForm

```
In[5]:= LDot[a, b]
```

```
Out[5]= a.b
```

```
In[4]:=  $\langle u[\mathbf{p}, \mathbf{m}], \gamma_\mu, v[\mathbf{p}, \mathbf{m}] \rangle$  // InputForm  
FermionLine[{1, p, m}, {-1, p, m},  
DiracMatrix[LTensor[DiracG,  $\mu$ ]]]
```

Traditional form also available!

```
DiracMatrix[ $\gamma_\mu$ , (k + p). $\gamma$ ,  $\gamma_\nu$ , PL] //  
TraditionalForm
```

$\gamma^\mu (\not{k} + \not{p}) \gamma^\nu P_L$
↑
 $\gamma \cdot p$

F) Integrated documentation

General Tensor Operations and Symbols

LTensor (\square), **LDot** (\square, \square), **LScalarQ** — Lorentz tensors and scalar products

Contract — contract tensors with repeated indices

Transverse, **Longitudinal** — project 2nd or higher rank Lorentz tensors

MandelstamRelations — express scalar products in terms of Mandelstam invariants

$g_{\square, \square}$ * $\epsilon_{\square, \square, \square, \square}$ — metric tensor, and Levi-Civita symbol

Computing One-Loop Integrals

LoopIntegrate, **LoopRefine** — routines for one loop integrals

LoopRefineSeries — compute series expansions of loop integrals

$d = \epsilon + \mu$ — dimensional regularization symbols

Dirac Algebra and Fermion Spinors

Spur — compute traces over product of Dirac matrices

Projector — project fermion self energy and vertex functions onto form factors

$\mathbf{1}$ * γ_\square * γ_5 * **PL** * **PR** * $\sigma_{\square, \square}$ — objects in spinor space

DiracMatrix, **FermionLine** (u, v), **FermionLineProduct**, **FermionLineExpand** — open fermion chains

Generated Symbols

PVA * **PVB** * **PVC** * **PVD** * **PVX** — symbolic Passarino-Veltman functions

Special Functions and Abbreviations »

Kallen λ * **DiscB** * **ScalarC0** * **ScalarD0** * ...

What is *Package-X* ?

skip

Selected features of *Package-X*

Package-X can generate analytic expressions for **arbitrarily high rank** tensor integrals **with up to four distinct propagators**, each with **arbitrary integer weight**, giving UV and IR divergent, and finite parts at **arbitrary (real-valued) kinematic points**, and can construct multivariable Taylor series expansions near **arbitrary (non-singular) kinematic point** to **arbitrary order**.

Package-X can calculate traces of products of gamma matrices, and perform tensor algebraic operations on open fermion lines.

Package-X can numerically evaluate scalar basis functions with either machine precision (fast) and arbitrary precision (slow) at any kinematic point.

What is *Package-X* ?

Applications: anything... except fully differential cross sections

- shift in vacuum expectation values
- shift in pole mass
- wavefunction renormalization
- electroweak oblique parameters
- particle electromagnetic moments (EDMs, polarizability, ...)
- 2 body decays, 3 body decays
- counterterms
- wilson coefficients
- cross sections at threshold
- ...

Motivation:

Until *Package-X*, there was no comprehensive one-loop package available.

There are semi-analytic packages (like FeynCalc,...) and numerical ones (like LoopTools,...)

- Sometimes difficult to use
- Does not always get an answer, like at vanishing Gram determinant, etc...

New in *Package-X* 2.1

A) Command-line readiness

Before, names of keywords were the symbols

```
 $\epsilon_{\mu,\nu,\rho,\sigma} g_{\rho,\nu} \text{DiracMatrix}[\gamma_\nu, \sigma_{\rho,\sigma}] // \text{FullForm}$   
  
Times[DiracMatrix[LTensor[\[Gamma], \[Nu]],  
  LTensor[\[Sigma], \[Rho], \[FinalSigma]],  
  LTensor[\[DoubleStruckG], \[Nu], \[Rho]],  
  LTensor[\[CurlyEpsilon], \[Mu],  
    \[Nu], \[Rho], \[FinalSigma]]]
```

*Now, keywords have proper full names,
and parse into full names by front-end.*

```
 $\epsilon_{\mu,\nu,\rho,\sigma} g_{\rho,\nu} \text{DiracMatrix}[\gamma_\nu, \sigma_{\rho,\sigma}] // \text{FullForm}$   
  
Times[DiracMatrix[LTensor[DiracG, \[Nu]],  
  LTensor[DiracS, \[Rho], \[FinalSigma]],  
  LTensor[MetricG, \[Nu], \[Rho]], LTensor[  
  LeviCivitaE, \[Mu], \[Nu], \[Rho], \[FinalSigma]]]
```

All old front-end notebooks continue to work,
but now also convenient to use command-line and edit .m (.wl) files:

▼ Square amplitude

```
applyPolarizationCompleteness[expr_, List[p_, 0, idx_]] :=  
  Replace[expr, {  
    HoldPattern[rest_. Power[LTensor[MetricG, PatternSequence[idx, a_]] | Pat  
    HoldPattern[rest_. LTensor[v_, idx]^2] :> (-LDot[v, v]) rest,  
    HoldPattern[rest_. LTensor[MetricG, PatternSequence[idx, a_]] | PatternSe  
    HoldPattern[rest_. LTensor[MetricG, PatternSequence[idx, a_]] | PatternSe  
    HoldPattern[rest_. LTensor[MetricG, idx, idx]] :> rest (-Dim),  
    HoldPattern[rest_. LTensor[v_, idx] * LTensor[w_, idx]] :> (-LDot[v, w])  
  
    expr :> ReplaceOnce[expr, LTensor[vec_, l____, idx, r____] :> -LTensor[vec, l,  
  ]];
```

```
Mathematica 10.3.1 for Mac OS X x86 (64-bit)  
Copyright 1988-2015 Wolfram Research, Inc.
```

```
In[1]:= <<X`  
Package-X v2.1.0 [developer version], by Hiren H. Patel
```

```
In[2]:= FermionLine[{-1, p, m}, {1, p, m}, DiracMatrix[LTensor[DiracG, mu]] * LTensor[q,  
mu]
```

```
Out[2]= < v[p, m], \gamma(mu), u[p, m] > q(mu)
```

```
In[3]:= %//Contract
```

```
Out[3]= < v[p, m], \gamma.q, u[p, m] >
```

```
In[4]:=
```

New in *Package-X* 2.1

B) Calculate the discontinuity across normal threshold cut for any one-loop integral:

Set option Part→*Discontinuity[s]* to *LoopRefine*

$$\int \frac{d^d k}{(2\pi)^d} \frac{k^\mu k^\nu}{[(k+p)^2 - m^2] k^2}$$

1. `LoopIntegrate[kμ kν, k, {k+p, m}, {k, 0}]`
`pμ pν PVB[0, 2, p.p, 0, m] + gμ,ν PVB[1, 0, p.p, 0, m]`

2. `% /. p.p → t`
`pμ pν PVB[0, 2, t, 0, m] + gμ,ν PVB[1, 0, t, 0, m]`

3. `LoopRefine[%, Part → Discontinuity[t]]`

$$\frac{2 i \pi (-m^2 + t)^3 \text{HeavisideTheta}[-m^2 + t] p_\mu p_\nu}{3 t^3} -$$

$$\frac{i \pi (-m^2 + t)^3 \text{HeavisideTheta}[-m^2 + t] g_{\mu,\nu}}{6 t^2}$$

Applications:

- Compute tree-level cross sections (optical theorem)
- Leptogenesis: only need imaginary parts of 1-loop diagrams

Important!

Discontinuity

`Discontinuity[s]`

is a setting for `Part` to `LoopRefine` to compute the discontinuity across the normal threshold cut of `s`.

`Discontinuity[s, t]`

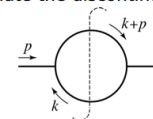
computes the Mandelstam double spectral function in overlapping channels `s` and `t`.

Details and Options

Examples (7)

Basic Examples (1)

Calculate the discontinuity across the normal threshold cut of the scalar two-point function



```
In[1]:= LoopIntegrate[1, k, {k+p, m}, {k, m}]
LoopRefine[%, Part → Discontinuity[p.p]]
```

```
Out[1]= PVB[0, 0, p.p, m, m]
```

COLLIER

April 2016, COLLIER released!



arXiv.org > hep-ph > arXiv:1604.06792

Search
(Help | Adv)

High Energy Physics - Phenomenology

Collier: a fortran-based Complex One-Loop Library in Extended Regularizations

Ansgar Denner, Stefan Dittmaier, Lars Hofer

- Numerical library of Passarino-Veltman functions (scalar *and* tensor integrals)
- Arbitrarily high rank *and* arbitrarily many legs
- Dimensional regularization *and* mass regularization
- Complex internal masses
- Dedicated expansions near vanishing Gram/Cayley determinants
- **Two** independent implementations (and nothing to do with FF library)

Bring to *Mathematica*! \implies CollierLink

CollierLink

Three features -

1. Direct numerical evaluation of Passarino-Veltman functions

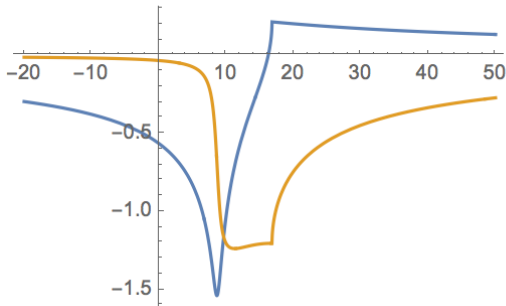
```
<< X`
<< CollierLink`
Package-X v2.1.0 [ALPHA], by Hiren H. Patel
For more information, see the guide
CollierLink v1.0.0 [ALPHA], by Hiren H. Patel
```

C_{0011}

```
PVC[1, 2, 0, 0, 0, 3.2, 1.0 - .2 I, 1.0 - .2 I, 2.0]
-0.0155049 + 0.010052 i
```

complex masses

```
Plot[{Re[PVC[0, 0, 0, 5.3, 12.1, s, 1.1, 1.2, 3.0]],
      Im[PVC[0, 0, 0, 5.3, 12.1, s, 1.1, 1.2, 3.0]]}, {s, -20, 50}]
```



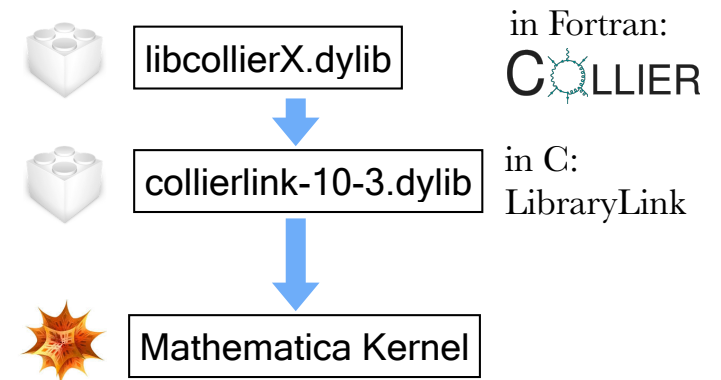
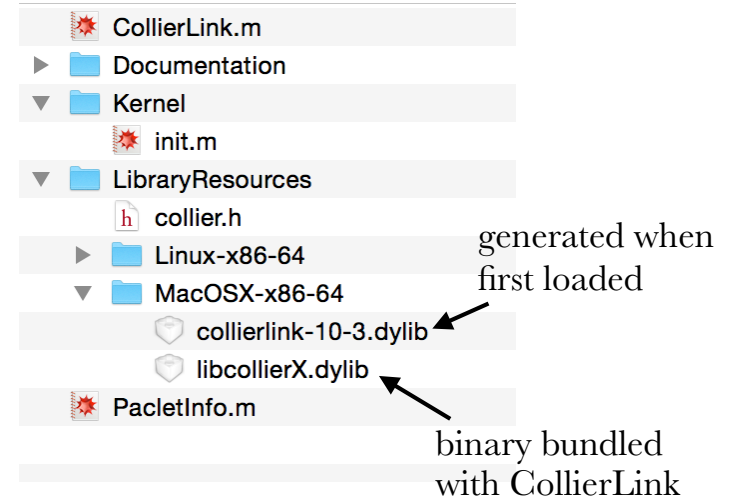
```
PVD[0, 0, 0, 0, 1, 2, 3, 4, 5, 6, 7, 8, 9, 10.0]
```

```
COLLIER::nocase : Error flag -10: Unable to
evaluate PVD[0, 0, 0, 0, 1., 2., 3., 4., 5., 6., 7., 8., 9., 10.];
case is not supported or implemented.
```

Indeterminate

Hiren Patel

under the hood...



CollierLink

Three features -

2. Automatic code generation, compilation, and relinking

$$\int \frac{d^d k}{(2\pi)^d} \frac{(k^2 + k.p + k.q + p.q)^2}{(k+p)^2 [(k+q)^2 - m^2] k^2}$$

```
myIntegral = LoopIntegrate[(k.k + k.p + k.q + p.q)^2, k, {k+p, 0}, {k+q, m}, {k, 0}] /.
```

```
{p.p -> m^2, q.q -> 0, p.q -> 1/2 (m^2 - t)}
```

```
3/4 PVA[0, 0] + 1/4 PVA[0, m] + (m^2/4 + 1/2 (m^2 - t)) PVB[0, 0, 0, 0, m] +
(-m^2/4 + 1/2 (m^2 - t)) PVB[0, 0, m^2, 0, 0] + 1/4 (m^2 - t) PVB[0, 1, 0, 0, m] +
1/4 (m^2 - t) PVB[0, 1, m^2, 0, 0] + 1/4 (m^2 - t)^2 PVC[0, 0, 0, m^2, t, 0, 0, 0, m]
```

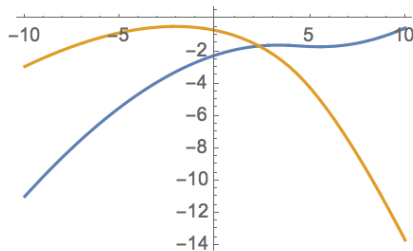
pretend
this is
many
pages
long...

```
f = CollierCompile[{t, m}, Evaluate[myIntegral]]
```

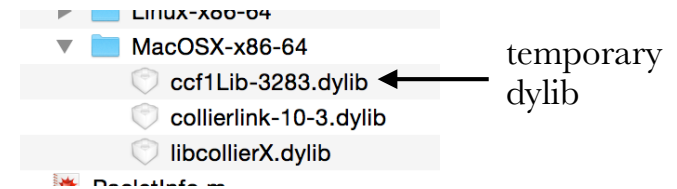
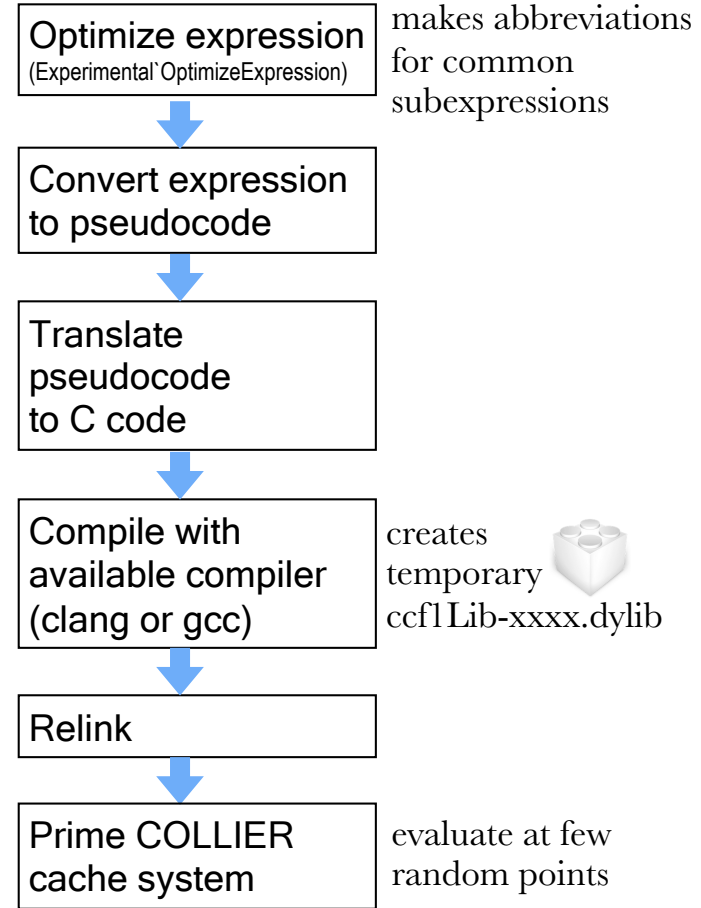
```
CollierCompiledFunction[ Variables (2): {t, m}
Denom: 3 (triangle) Rank: 1 ]
```

Evaluates with machine-code speed!

```
Plot[{Re[f[t, 2.1]], Im[f[t, 2.1]]}, {t, -10, 10}]
```



under the hood...



CollierLink


Three features -

2. *Automatic code generation, compilation, and relinking*


uses COLLIER's cache system —

Each CollierCompiledFunction gets its own cache


```
CollierCompile[{s, m}, PVA[0, m]]
```

```
CollierCompiledFunction [ -  Variables (2): {s, m}  
Denom: 1 (tadpole) Rank: 0  
TN_cll calls: 1 cache_no: 1  
Code lines: 2 File size: 14.80 kB
```

```
CollierCompile[{s, m}, PVC[1, 0, 0, 0, 0, s, m, m, m]]
```

```
CollierCompiledFunction [ -  Variables (2): {s, m}  
Denom: 3 (triangle) Rank: 2  
TN_cll calls: 1 cache_no: 2  
Code lines: 2 File size: 14.84 kB
```

```
CollierCompile[{s, m}, PVB[0, 0, s, m, m]]
```

```
CollierCompiledFunction [ -  Variables (2): {s, m}  
Denom: 2 (bubble) Rank: 0  
TN_cll calls: 1 cache_no: 3  
Code lines: 2 File size: 14.84 kB
```

No conflicts

And CollierCompiledFunction can be parallelized!

CollierCompiledFunction

```
CollierCompiledFunction[args, ...]
```

represents compiled code for evaluating a compiled function using the COLLIER library.

Compile the infrared divergent scalar triangle function:

```
In[1]:= myfun = CollierCompile[{qSq}, PVC[0, 0, 0, 0, 0, qSq, 0, 0, 0]]
```

```
Out[1]= CollierCompiledFunction [ +  Variables (1): {qSq}  
Denom: 3 (triangle) Rank: 0
```

Load CollierLink on all parallel kernels with `ParallelNeeds`, and distribute the definition of `myFun` to parallel kernels:

```
In[2]:= ParallelNeeds["CollierLink`"]  
DistributeDefinitions[myfun]
```

```
Out[2]= {myfun}
```

Compute a table of values in parallel:

```
In[3]:= ParallelTable[myfun[qSq], {qSq, 5.1, 10.3}]
```

```
Out[3]= {-0.868639 - 1.00361 i, -0.675789 - 0.931296 i, -0.540322 - 0.867299 i,  
-0.440657 - 0.811331 i, -0.364729 - 0.762362 i, -0.305283 - 0.719311 i}
```

Each parallel kernel gets its own copy of dylib

CollierLink

Three features -

3. Automatic Fortran code generation

CollierCodeGenerate [{t, m}, Evaluate[myIntegral]]

```
!!*****!!
!! This code was automatically created using Package-X v2.1.0 CollierLink !!
!! routine CollierCodeGenerate[] by Hiren H. Patel, and requires the !!
!! Fortran library COLLIER v1.1 by A. Denner, S. Dittmaier, L. Hofer !!
!! for numerical evaluation. !!
!!-----!!
!! Language: Fortran Creation date: Aug-21-2017 21:15 !!
!! Function name: generatedFunction Arguments (2): {t,m} !!
!! Author: hhpate1 !!
!! Initialization requirement: Init_c11(3,1,') !!
!!*****!!

DOUBLE COMPLEX FUNCTION generatedFunction(inputVar1, inputVar2)

USE COLLIER
IMPLICIT NONE

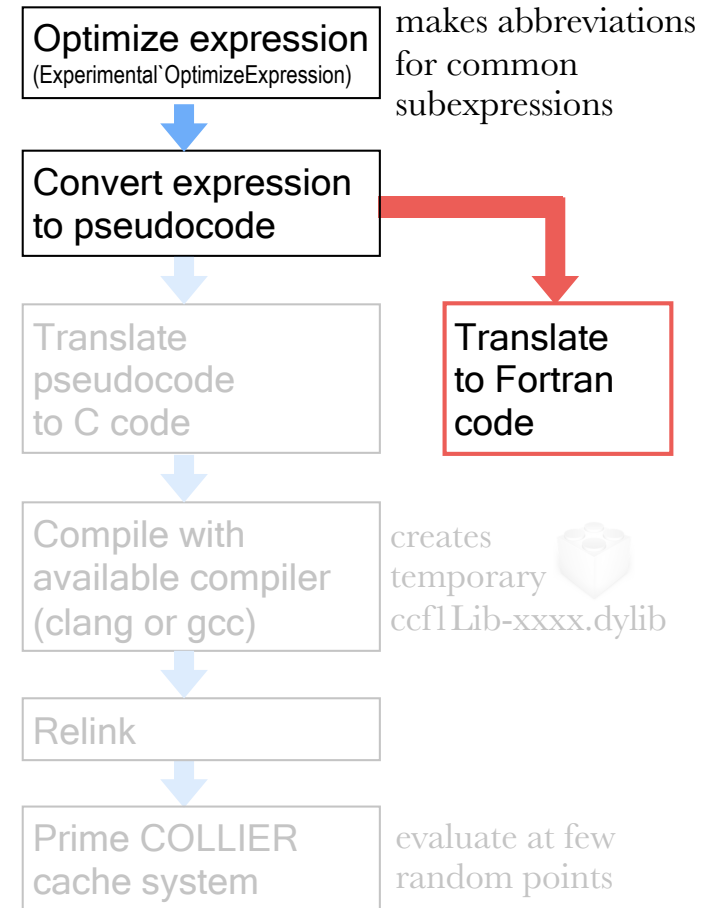
DOUBLE COMPLEX, INTENT(IN) :: inputVar1 ! t
DOUBLE COMPLEX, INTENT(IN) :: inputVar2 ! m

!Allocate memory for calculated results of Passarino-Veltman,
! coefficient functions, and their UV divergent parts
DOUBLE COMPLEX :: pvx1(1), pvx1uv(1) ! PVA[_ ,0]
DOUBLE COMPLEX :: pvx2(1), pvx2uv(1) ! PVA[_ ,m]
DOUBLE COMPLEX :: pvx3(2), pvx3uv(2) ! PVB[_ ,_,0,0,m]
DOUBLE COMPLEX :: pvx4(2), pvx4uv(2) ! PVB[_ ,_,Compile`optVar5,0,0]
DOUBLE COMPLEX :: pvx5(3), pvx5uv(3) ! PVC[_ ,_,_,Compile`optVar5,t,0,0,0,m]

!Local variables
DOUBLE COMPLEX :: localVar1 ! optVar5
DOUBLE COMPLEX :: localVar2 ! optVar7
DOUBLE COMPLEX :: localVar3 ! optVar8
DOUBLE COMPLEX :: localVar4 ! optVar9
DOUBLE COMPLEX :: localVar5 ! optVar21

!Evaluation code
localVar1 = (inputVar2**2)
localVar2 = (-inputVar1)
localVar3 = (localVar1+localVar2)
localVar4 = ((1.d0/2.d0)*localVar3)
localVar5 = (localVar3**2)
```

under the hood...



Idea is to be able to easily plug this code into popular event generators MCFM, MadGraph, ...

Summary

- ★ Introduced *Package-X*, described its salient features
- ★ Described updates in upcoming release *Package-X* v2.1
- ★ Introduced new package, CollierLink
 1. *Direct numerical evaluation of Passarino-Veltman functions*
 2. *Automatic code generation, compilation, and relinking*
 3. *Automatic Fortran code generation*

Conclusions:

Together with CollierLink, *Package-X* provides a user-friendly yet *truly comprehensive tool* for the calculation of one loop integrals.

Applications: anything... ~~except fully differential cross sections~~