RB-SFA: High Harmonic Generation in the Strong Field Approximation via *Mathematica*

Introduction

Readme

RB-SFA is a compact and flexible *Mathematica* package for calculating High Harmonic Generation emission within the Strong Field Approximation. It combines *Mathematica*'s analytical integration capabilities with its numerical calculation capacities to offer a fast and user-friendly plug-and-play solver for calculating harmonic spectra and other properties. In addition, it can calculate first-order nondipole corrections to the SFA results to evaluate the effect of the driving laser's magnetic field on harmonic spectra.

The name RB-SFA comes from its first application (as Rotating Bicircular High Harmonic Generation in the Strong field Approximation) but the code is general so RB-SFA just stands for itself now. This first application was used to calculate the polarization properties of the harmonics produced by multi-colour circularly polarized fields, as reported in the paper

Spin conservation in high-order-harmonic generation using bicircular fields. E. Pisanty, S. Sukiasyan and M. Ivanov. *Phys. Rev. A* **90**, 043829 (2014), arXiv:1404.6242.

This code is dual-licensed under the GPL and CC-BY-SA licenses. If you use this code or its results in an academic publication, please cite the paper above or the GitHub repository where the latest version will always be available. An example citation is

E. Pisanty. RB-SFA: High Harmonic Generation in the Strong Field Approximation via *Mathematica*. https://github.com/episanty/RB-SFA (2016).

This software consists of this notebook, which contains the code and its documentation, a corresponding auto-generated package file. This notebook also contains a Usage and Examples section which explains how to use the code and documents the calculations used in the original publication.

Implementation

Supporting functions

Initialization

```
BeginPackage["RBSFA`"];
```

Version number

The command RBSFAversion prints the version of the RB-SFA package currently loaded and its timestamp

```
$RBSFAversion::usage = "$RBSFAversion prints the current
        version of the RB-SFA package in use and its timestamp  .";
$RBSFAtimestamp  ::usage = "$RBSFAtimestamp   prints the timestamp
        of the current version of the RB-SFA package.";
Begin["`Private`"];
$RBSFAversion:= "RB-SFA v2.1.3, " <> $RBSFAtimestamp  ;
End[];
```

Old syntax (in functional form RBSFAversion[]), deprecated

```
RBSFAversion::usage = "RBSFAversion[] has been deprecated in favour of $RBSFAversion.";
RBSFAversion::dprc = "RBSFAversion[] has been deprecated in favour of $RBSFAversion.";
Begin["`Private`"];
RBSFAversion[] := (Message[RBSFAversion::dprc]; $RBSFAversion);
End[];
```

The timestamp is updated every time the notebook is saved via an appropriate notebook option, which is set by the code below.

```
SetOptions[
    EvaluationNotebook[],
    NotebookEventActions→{{"MenuCommand  ", "Save"}:→(
            NotebookWrite[
              Cells[CellTags→"version-timestamp  "][[1]],
              Cell[
                BoxData[RowBox[{"Begin[\"`Private`\"];\n$RBSFAtimestamp  =\""<>
                        DateString[]<>"\";\nEnd[];"}]]
                , "Input", InitializationCell→True, CellTags→"version-timestamp  "
              ], None, AutoScroll→False];
            NotebookSave[]
          ), PassEventsDown→True}
  ];
```

To reset this behaviour to normal, evaluate the cell below

```
SetOptions[EvaluationNotebook[],
  NotebookEventActions→{{"MenuCommand  ", "Save"}:→(NotebookSave[]), PassEventsDown→True}]
```

Timestamp

```
Begin["`Private`"];
$RBSFAtimestamp  = "Wed 25 Jan 2017 18:40:33";
End[];
```

Directory

```
$RBSFAdirectory::usage = "$RBSFAdirectory is the directory
        where the current RB-SFA package instance is located.";
```

```
Begin["`Private`"];
With[{softLinkTestString= StringSplit[StringJoin[ReadList[
               "! ls -la "<>StringReplace[$InputFileName , {" " → "\\ "}], String]], " -> "]},
    If[Length[softLinkTestString] > 1, (*Testing in case $InputFileName
       is a soft link to the actual directory.*)
      $RBSFAdirectory= StringReplace[DirectoryName [softLinkTestString[[2]]], {" " → "\\ "}],
      $RBSFAdirectory= StringReplace[DirectoryName [$InputFileName ], {" " → "\\ "}];
    ]];
End[];
```

Git commit hash and message

```
$RBSFAcommit  ::usage = "$RBSFAcommit   returns the git commit
        log at the location of the RB-SFA package if there is one.";
$RBSFAcommit  ::OS = "$RBSFAcommit   has only been tested on Linux.";

Begin["`Private`"];
$RBSFAcommit  := (If[$OperatingSystem ≠ "Unix", Message[$RBSFAcommit  ::OS]];
      StringJoin[
        Riffle[ReadList["!cd "<>$RBSFAdirectory<>" && git log -1", String], {"\n"}]]);
End[];
```

Standard function (re)definitions

ConstantArray

This redefines ConstantArray to take the corner case of an empty dimensions list, which returns an error code (and an unevaluated ConstantArray) for Mathematica versions under 10.1.0 (cf. mma.se/q/133078).

```
Quiet[Check[
      ConstantArray[0, {}];,
      Unprotect[ConstantArray];
      ConstantArray[Private`x_, {}] := Private`x;
      Protect[ConstantArray];
    ]];
```

Similarly, this needs to be put inside an initialization code for any parallelized subkernels that may get launched later (cf. mm.se/q/131856).

```
Parallelize;
Parallel`Developer`$InitCode= Hold[
      Quiet[Check[
            ConstantArray[0, {}];,
            Unprotect[ConstantArray];
            ConstantArray[Private`x_, {}] := Private`x;
            Protect[ConstantArray];
          ]];
    ];
```

ReIm

This adds the definition of ReIm for those versions (<10.1) that don't have it.

```
If[
  Context[ReIm ] =!= "System` ",
  ReIm ::usage =
    "\!\(\*RowBox[{\"ReIm \", \"[\", StyleBox[\"z\", \"TI\"], \"]\"}]\) gives the list
        \!\(\*RowBox[{\"{\", RowBox[{RowBox[{\"Re\", \"[\", StyleBox[\"z\",
        \"TI\"], \"]\"}], \",\", RowBox[{\"Im \", \"[\", StyleBox[\"z\", \"TI\"],
        \"]\"}]}], \"}\"}]\) of the number  \!\(\*StyleBox[\"z\", \"TI\"]\).";
  ReIm [Private`z_] := {Re[Private`z], Im [Private`z]};
  SetAttributes[ReIm , Listable];
  Protect[ReIm ];
]
```

AssociationTranspose

```
ClearAll[AssociationTranspose]

AssociationTranspose::usage =
    "AssociationTranspose[association] transposes the given two-level
        association of associations.";
AssociationTranspose::wrngshp = "Input `1` is the wrong shape; it must  be
        an association all of whose Values are valid associations.";
Begin["`Private`"];
AssociationTranspose[association_?(
          And@@(AssociationQ/@Join[{#}, Values[#]]) &
        )] := GroupBy[
        Join@@Thread/@Normal //@association,
        {First@*Last, First}
      ][[All, All, 1, 2, 2]];
AssociationTranspose[association_] :=
  "Doesn't display; cf. mm  .se/q/29321 for details"/;
    Message[AssociationTranspose::wrngshp, association]
End[];
```

The above function is taken from http://mathematica.stackexchange.com/a/86526 by http://mathematica.stackexchange.com/users/121/mr-wizard. *Mathematica* 10.1 and higher, it can be replaced by a Query[Transpose] construct as below, but *Mathematica* 10.0, despite having most of the Association code, is unable to transpose ragged associations using that construct.

```
(*
AssociationTranspose[association_]:=DeleteMissing[
    Query[Transpose][
      association
    ]
    ,2]
*)
```

KeyValueMap

This exists in *Mathematica* 10.1 and later, but it's nice to have it on version 10.0 so this is a back-port for versions that do not have it.

```
If[
  $VersionNumber < 10.1,
  KeyValueMap::usage =
    "\!\(\*RowBox[{\"KeyValueMap\", \"[\", RowBox[{StyleBox[\"f\",\"TI\"], \",\",
        RowBox[{\"<|\",RowBox[{RowBox[{SubscriptBox[StyleBox[\"key\", \"TI\"],
        StyleBox[\"1\", \"TR\"]], \"→\", SubscriptBox[StyleBox[\"val\",
        \"TI\"], StyleBox[\"1\", \"TR\"]]}], \",\",
        RowBox[{SubscriptBox[StyleBox[\"key\", \"TI\"], StyleBox[\"2\",
        \"TR\"]], \"→\", SubscriptBox[StyleBox[\"val\", \"TI\"], StyleBox[\"2\",
        \"TR\"]]}], \",\", StyleBox[\"...\", \"TR\"]}], \"|>\"}]}], \"]\"}]\)
        gives the list \!\(\*RowBox[{\"{\", RowBox[{RowBox[{StyleBox[\"f\",
        \"TI\"], \"[\", RowBox[{SubscriptBox[StyleBox[\"key\", \"TI\"],
        StyleBox[\"1\", \"TR\"]], \",\", SubscriptBox[StyleBox[\"val\", \"TI\"],
        StyleBox[\"1\", \"TR\"]]}], \"]\"}], \",\", RowBox[{StyleBox[\"f\",
        \"TI\"], \"[\", RowBox[{SubscriptBox[StyleBox[\"key\", \"TI\"],
        StyleBox[\"2\", \"TR\"]], \",\", SubscriptBox[StyleBox[\"val\",
        \"TI\"], StyleBox[\"2\", \"TR\"]]}], \"]\"}], \",\", StyleBox[\"...\",
        \"TR\"]}], \"}\"}]\). (Note: function backported from  v10.1+.)

\!\(\*RowBox[{\"KeyValueMap\", \"[\", StyleBox[\"f\", \"TI\"], \"]\"}]\) represents an
        operator form  of KeyValueMap that can be applied to an expression.";
  KeyValueMap::invak = "The argument  `1` is not a valid association";

  Begin["`Private`"];
  KeyValueMap[f_, assoc_?AssociationQ] := Map[Apply[f], Normal [assoc]];
  KeyValueMap[f_][assoc_?AssociationQ] := KeyValueMap[f, assoc];
  KeyValueMap[f_, assoc__] := "Doesn't display; cf. mm  .se/q/29321 for details" /;
      Message[KeyValueMap::invak, assoc];
  End[];
]
```

Dipole transition matrix elements

Default DTME, for a hydrogenic 1s state

```
hydrogenicDTME::usage =
    "hydrogenicDTME[p,κ] returns the dipole transition matrix element for
        a 1s hydrogenic state of ionizationpotential Ip=1/2 κ^2.

hydrogenicDTME[p,κ,{n,l,m }] returns the dipole transition matrix element for an
        n,l,m hydrogenic state of ground-state ionizationpotential Ip=1/2 κ^2.

hydrogenicDTME[p,κ,n,l,m ] returns the dipole transition matrix element for an
        n,l,m hydrogenic state of ground-state ionizationpotential Ip=1/2 κ^2.";
hydrogenicDTMERegularized::usage = "hydrogenicDTMERegularized[p,κ] returns the
        dipole transition matrix element for a 1s hydrogenic state of
        ionizationpotential Ip=1/2 κ^2, regularized to remove the denominator
        of 1/(p^2+κ^2)^3, where the saddle-point solutions are singular.

hydrogenicDTMERegularized[p,κ,{n,l,m }] returns the dipole transition matrix element for
        an n,l,m hydrogenic state of ground-state ionizationpotential Ip=1/2 κ^2,
        regularized to remove factors of (p^2+κ^2) from the denominator .

hydrogenicDTMERegularized[p,κ,n,l,m ] returns the dipole transition matrix element for an
        n,l,m hydrogenic state of ground-state ionizationpotential Ip=1/2 κ^2,
        regularized to remove factors of (p^2+κ^2) from the denominator .";
Begin["`Private`"];
hydrogenicDTME[p_List, κ_] := 8 i/π √(2 κ^5) p / (Total[p^2] + κ^2)^3
hydrogenicDTME[p_?NumberQ , κ_] := 8 i/π √(2 κ^5) p / (p^2 + κ^2)^3
hydrogenicDTMERegularized[p_List, κ_] := 8 i/π √(2 κ^5) p / 1
hydrogenicDTMERegularized[p_?NumberQ , κ_] := 8 i/π √(2 κ^5) p / 1
End[];
```

For a gaussian orbital

```
gaussianDTME::usage =
    "gaussianDTME[p,κ] returns the dipole transition matrix element for a gaussian
        state of characteristic size 1/κ.";
Begin["`Private`"];
gaussianDTME[p_List, κ_] := -i (4 π)^(3/4) κ^(-7/2) p Exp[-Total[p^2]/(2 κ^2)]
gaussianDTME[p_?NumberQ , κ_] := -i (4 π)^(3/4) κ^(-7/2) p Exp[-p^2/(2 κ^2)]
End[];
```

SolidHarmonicS

This function implements the solid harmonic $S_{l,m}(\boldsymbol{r}) = r^l Y_{l,m}(\theta, \phi)$, which is a homogeneous polynomial of degree l, and lends itself much better to symbolic differentiation than explicit spherical harmonics.

Code provided by J.M. at http://mathematica.stackexchange.com/a/124336/1000 under the WTFPL.

```
SolidHarmonicS::usage =
    "SolidHarmonicS[l,m,x,y,z] calculates the solid harmonic S_lm(x,y,z)=r^l Y_lm(x,y,z).

SolidHarmonicS[l,m,{x,y,z}] does the same .";
Begin["`Private`"];
SolidHarmonicS[λ_Integer, μ_Integer, x_, y_, z_] /; λ ≥ Abs[μ] :=
  Sqrt[(2 λ+1)/(4 π)] Sqrt[Gamma[λ-Abs[μ]+1]/Gamma[λ+Abs[μ]+1]] 2^-λ (-1)^((μ-Abs[μ])/2) ×
    If[Rationalize[μ] == 0, 1, (x+Sign[μ] i y)^Abs[μ]] ×
    Sum[
      (-1)^(μ+k) Binomial[λ, k] Binomial[2 λ-2 k, λ] Pochhammer[λ-Abs[μ]-2 k+1, Abs[μ]] ×
        If[Rationalize[k] == 0, 1, (x^2+y^2+z^2)^k] ×
        If[Rationalize[λ-Abs[μ]-2 k] == 0, 1, z^(λ-Abs[μ]-2 k)]
      , {k, 0, Quotient[λ, 2]}]
SolidHarmonicS[λ_Integer, μ_Integer, {x_, y_, z_}] /; λ ≥ Abs[μ] :=
  SolidHarmonicS[λ, μ, x, y, z]
End[];
```

hydrogenicΨ and hydrogenicY (momentum-space wavefunctions)

This implements the dipole transition matrix element from an arbitrary hydrogenic orbital n, l, m , where the ground-state ionization potential is given by $I_p = \frac{1}{2}\kappa^2$, as described in Luke Chipperfield's PhD thesis (Imperial College London, 2008, p. 52). This code uses partial memoization as in mm.se/q/21782.

```
hydrogenicΨ::usage =
    "hydrogenicΨ[n,l,m ,κ,px,py,pz] calculates the momentum   -space wavefunction
        Ψ(p)=⟨p|nlm ⟩ for a hydrogenic atom  with ionizationpotential κ^2/2.

hydrogenicΨ[n,l,m ,κ,{px,py,pz}] calculates the momentum   -space wavefunction
        Ψ(p)=⟨p|nlm ⟩ for a hydrogenic atom  with ionizationpotential κ^2/2.";
Begin["`Private`"];
hydrogenicΨ[n_, l_, m_ , κκ_, ppx_, ppy_, ppz_] := Block[{κ, px, py, pz},
      hydrogenicΨ[n, l, m , κ_, px_, py_, pz_] = Simplify [
          -SolidHarmonicS [l, m , px, py, pz] (-i)^l π 2^(2 l+4) l! / (2 π κ)^(3/2) Sqrt[n (n-l-1) ! / (n+l) !]
            κ^(l+4) / (px^2 + py^2 + pz^2 + κ^2)^(l+2) GegenbauerC[n-l-1, l+1, (px^2 + py^2 + pz^2 - κ^2)/(px^2 + py^2 + pz^2 + κ^2)]
        ];
      hydrogenicΨ[n, l, m , κκ, ppx, ppy, ppz]
    ];
hydrogenicΨ[n_, l_, m_ , κ_, {px_, py_, pz_}] := hydrogenicΨ[n, l, m , κ, px, py, pz];
End[];
```

Regularized version, removing the powers of $p^2 + \kappa^2$ in the denominator, to eliminate poles at the saddle-point momentum $p = i\,\kappa$.

```
hydrogenicΨRegularized::usage =
    "hydrogenicΥRegularized[n,l,m ,κ,px,py,pz] calculates the momentum   -space
        wavefunction Ψ(p)=⟨p|nlm ⟩ for a hydrogenic atom  with
        ionizationpotential κ^2/2, multiplied by (p^2+κ^2)^(n+1)
        to remove  any factors of p^2+κ^2 in the denominator .

hydrogenicΥRegularized[n,l,m ,κ,{px,py,pz}] calculates the momentum   -space wavefunction
        Ψ(p)=⟨p|nlm ⟩ for a hydrogenic atom  with ionizationpotential κ^2/2, multiplied
        by (p^2+κ^2)^(n+1) to remove  any factors of p^2+κ^2 in the denominator .";
Begin["`Private`"];
hydrogenicΨRegularized[n_, l_, m_ , κκ_, ppx_, ppy_, ppz_] := Block[{κ, px, py, pz},
      hydrogenicΨRegularized[n, l, m , κ_, px_, py_, pz_] = Simplify [Cancel[
          -SolidHarmonicS [l, m , px, py, pz] (-i)^l π 2^(2 l+4) l! / (2 π κ)^(3/2) Sqrt[n (n-l-1) ! / (n+l) !]
            κ^(l+4) (px^2 + py^2 + pz^2 + κ^2)^(n-l-1) GegenbauerC[n-l-1, l+1, (px^2 + py^2 + pz^2 - κ^2)/(px^2 + py^2 + pz^2 + κ^2)]
        ]];
      hydrogenicΨRegularized[n, l, m , κκ, ppx, ppy, ppz]
    ];
hydrogenicΨRegularized[n_, l_, m_ , κ_, {px_, py_, pz_}] :=
    hydrogenicΨRegularized[n, l, m , κ, px, py, pz];
End[];
```

Upsilon function, given by $\Upsilon(\boldsymbol{p}) = \left(\frac{1}{2}\boldsymbol{p}^2 + I_p\right)\Psi(\boldsymbol{p}) = \frac{1}{2}\left(\boldsymbol{p}^2 + \kappa^2\right)\langle \boldsymbol{p} \mid n, l, m \rangle$, which can be used in the form $\Upsilon(\boldsymbol{p} + \boldsymbol{A}(t'))$ as a replacement for the ionization dipole $\boldsymbol{d}(\boldsymbol{p} + \boldsymbol{A}(t'))\cdot\boldsymbol{F}(t')$, particularly for cases where the latter is singular but the former is not. (For details cf. arXiv:1304.2413, appendix A.)

```
hydrogenicΥ::usage =
    "hydrogenicΥ[n,l,m ,κ,px,py,pz] calculates the Upsilon function Υ(p)=(1/2 p^2+I_p)⟨p|nlm ⟩
        for a hydrogenic atom  with ionizationpotential κ^2/2.

hydrogenicΥ[n,l,m ,κ,{px,py,pz}] calculates the Upsilon function Υ(p)=(1/2 p^2+I_p)⟨p|nlm ⟩
        for a hydrogenic atom  with ionizationpotential κ^2/2.";
Begin["`Private`"];
hydrogenicΥ[n_, l_, m_ , κ_, px_, py_, pz_] :=
    1/2 (px^2 + py^2 + pz^2 + κ^2) hydrogenicΨ[n, l, m , κ, px, py, pz];
hydrogenicΥ[n_, l_, m_ , κ_, {px_, py_, pz_}] := hydrogenicΥ[n, l, m , κ, px, py, pz];
End[];
```

hydrogenicDTME for arbitrary states

```
Begin["`Private`"];
hydrogenicDTME[{ppx_, ppy_, ppz_}, κκ_, n_, l_, m_ ] := Block[{κ, px, py, pz},
      hydrogenicDTME[{px_, py_, pz_}, κ_, n, l, m ] =
        Simplify [Grad[hydrogenicΥ[n, l, m , κ, px, py, pz], {px, py, pz}]];
      hydrogenicDTME[{ppx, ppy, ppz}, κκ, n, l, m ]
    ];
hydrogenicDTME[{px_, py_, pz_}, κ_, {n_, l_, m_ }] := hydrogenicDTME[{px, py, pz}, κ, n, l, m ];
End[];
```

Regularized version, removing the powers of $p^2 + \kappa^2$ in the denominator, to eliminate poles at the saddle-point momentum $p = i\,\kappa$.

```
Begin["`Private`"];
hydrogenicDTMERegularized[{px_, py_, pz_}, κ_, n_, l_, m_ ] :=
    (px^2 + py^2 + pz^2 + κ^2)^(n+1) hydrogenicDTME[{px, py, pz}, κ, n, l, m ];
hydrogenicDTMERegularized[{px_, py_, pz_}, κ_, {n_, l_, m_ }] :=
    hydrogenicDTMERegularized[{px, py, pz}, κ, n, l, m ];
End[];
```

Various field envelopes

flatTopEnvelope

```
flatTopEnvelope::usage =
    "flatTopEnvelope[ω,num ,nRamp ] returns a Function object representing
        a flat-top envelope at carrier frequency ω lasting a total
        of num  cycles and with linear ramps  nRamp  cycles long.";
Begin["`Private`"];
flatTopEnvelope[ω_, num_ , nRamp_ ] := Function[t,
    Piecewise[{{0, t < 0}, {Sin[ω t/(4 nRamp)]^2, 0 ≤ t < 2π/ω nRamp }, {1, 2π/ω nRamp  ≤ t < 2π/ω (num  - nRamp )},
        {Sin[ω (2π/ω num  - t)/(4 nRamp)]^2, 2π/ω (num  - nRamp ) ≤ t < 2π/ω num }, {0, 2π/ω num  ≤ t}}]]
End[];
```

cosPowerFlatTop

```
cosPowerFlatTop::usage =
    "cosPowerFlatTop[ω,num ,power] returns a Function object representing
        a smooth  flat-top envelope of the form  1-Cos(ω t/2 num )^power";
Begin["`Private`"];
cosPowerFlatTop[ω_, num_ , power_] := Function[t, 1 - Cos[ω t / (2 num)]^power]
End[];
```

Field duration standard options

The standard options for the duration of the pulse and the resolution are

```
PointsPerCycle::usage =
    "PointsPerCycle is a sampling  option which specifies the number  of sampling
        points per cycle to be used in integrations.";
TotalCycles::usage = "TotalCycles is a sampling  option which specifies
        the total number  of periods to be integrated over.";
CarrierFrequency::usage = "CarrierFrequency is a sampling  option which
        specifies the carrier frequency to be used.";
Protect[PointsPerCycle, TotalCycles, CarrierFrequency];

standardOptions= {PointsPerCycle→ 90, TotalCycles → 1,
      CarrierFrequency→ 0.057, IntegrationPointsPerCycle→ Automatic };
```

PointsPerCycle dictates how many sampling points are used per laser cycle (at frequency CarrierFrequency, of the infrared laser), and it should be at least twice the highest harmonic of interest. The total duration is TotalCycles cycles. CarrierFrequency is the frequency of the fundamental laser, in atomic units.

harmonicOrderAxis

harmonicOrderAxis produces a list that can be used as a harmonic order axis for the given pulse parameters.

The length can be fine-tuned (to match exactly a spectrum, for instance, and get a matrix of the correct shape) using the correction option, or a TargetLength can be directly specified.

```
harmonicOrderAxis::usage =
    "harmonicOrderAxis[opt→value] returns a list of frequencies which can be used as
        a frequency axis for Fourier transforms , scaled in units of harmonic
        order, for the provided field duration and sampling  options.";
TargetLength::usage = "TargetLength is an option for harmonicOrderAxis  which
        specifies the total length required of the resulting list.";
LengthCorrection::usage = "LengthCorrection is an option for harmonicOrderAxis  which
        allows for manual  correction of the length of the resulting list.";
Protect[LengthCorrection, TargetLength];
Begin["`Private`"];
Options[harmonicOrderAxis] =
    standardOptions~Join~{TargetLength→Automatic , LengthCorrection→1};
harmonicOrderAxis::target =
    "Invalid TargetLength option `1`. This must  be a positive integer or Automatic .";
harmonicOrderAxis[OptionsPattern[]] :=
  Module[{num = OptionValue[TotalCycles], npp = OptionValue[PointsPerCycle]},
    Piecewise[{
        {1/num Range[0., Round[(npp num + 1)/2.] - 1 + OptionValue[LengthCorrection]],
          OptionValue[TargetLength] === Automatic},
        {Round[(npp num + 1)/2.]/num Range[0, OptionValue[TargetLength] - 1]/OptionValue[TargetLength],
          IntegerQ[OptionValue[TargetLength]] && OptionValue[TargetLength] ≥ 0}
      },
      Message[harmonicOrderAxis::target, OptionValue["TargetLength"]];
      Abort[]
    ]
  ]
End[];
```

frequencyAxis

frequencyAxis produces a list that can be used as a harmonic order axis for the given pulse parameters. Identical to harmonicOrderAxis but produces a frequency axis (in atomic units) instead.

```
frequencyAxis::usage =
    "frequencyAxis[opt→value] returns a list of frequencies which can be used
        as a frequency axis for Fourier transforms , in atomic  units of
        frequency, for the provided field duration and sampling  options.";
Begin["`Private`"];
Options[frequencyAxis] = Options[harmonicOrderAxis];
frequencyAxis[options:OptionsPattern[]] :=
  OptionValue[CarrierFrequency] harmonicOrderAxis[options]
End[];
```

timeAxis

timeAxis produces a list that can be used as a time axis for the given pulse parameters.

```
Quit
```

```
timeAxis::usage =
    "timeAxis[opt→value] returns a list of times which can be used as a time axis ";
TimeScale::usage = "TimeScale is an option for timeAxis which specifies the units the
        list should use: AtomicUnits by default, or LaserPeriods if required.";
AtomicUnits::usage = "AtomicUnits is a value for the option TimeScale of timeAxis
        which specifies that the times should be in atomic units of time .";
LaserPeriods::usage = "LaserPeriods is a value for the option TimeScale of timeAxis which
        specifies that the times should be in multiples of the carrier laser period.";
Protect[TimeScale, AtomicUnits, LaserPeriods];
Begin["`Private`"];
Options[timeAxis] =
    standardOptions~Join~{TimeScale → AtomicUnits, PointNumberCorrection → 0};
timeAxis::scale =
    "Invalid TimeScale option `1`. Available values are AtomicUnits and LaserPeriods";
timeAxis[OptionsPattern[]] := Block[{T = 2π/ω, ω = OptionValue[CarrierFrequency],
      num = OptionValue[TotalCycles], npp = OptionValue[PointsPerCycle]},
    Piecewise[{
          {1, OptionValue[TimeScale] === AtomicUnits},
          {1/T, OptionValue[TimeScale] === LaserPeriods}
        },
        Message[timeAxis::scale, OptionValue[TimeScale]]; Abort[]
      ] × Table[t
        , {t, 0, num 2π/ω, num/(num × npp + OptionValue[PointNumberCorrection]) 2π/ω}
      ]
  ]
End[];

tInit = 0;
tFinal = 2π/ω num ;
δt = (tFinal - tInit)/(num × npp + OptionValue[PointNumberCorrection]); (*integration and looping timestep *)
```

getSpectrum

getSpectrum takes a time-dependent dipole list and returns its Fourier transform in absolute-value-squared. It takes as options

- pulse parameters ω, TotalCycles and PointsPerCycle,
- a polarization parameter ϵ, which gives an unpolarized spectrum when given False, or polarizes along an ellipticity vector ϵ (this is meant primarily to select right- and left-circularly polarized spectra using $\epsilon = \{1, i\}$ and $\epsilon = \{1, -i\}$ respectively),
- a DifferentiationOrder, which can return the dipole value (default, = 0), velocity (= 1), or acceleration (= 2),
- a power of ω, ωPower, with which to multiply the spectrum before returning it (which should be equivalent to DifferentiationOrder except for pathological cases), and
- a ComplexPart function to apply immediately after differentiation (default is the identity function, but Re, Im, or Abs[#]^2 & are reasonable choices).

If no option is passed to ωPower and DifferentiationOrder, the pulse parameters do not really affect the output, except by a global factor of TotalCycles.

```
getSpectrum::usage = "getSpectrum[DipoleList] returns the power spectrum of DipoleList.";
Polarization::usage =
```

```
    "Polarization is an option for getSpectrum  which specifies a polarization
        vector along which to polarize the dipole list. The default,
        Polarization→False, specifies an unpolarized spectrum .";
ComplexPart ::usage = "ComplexPart  is an option for getSpectrum  which specifies
        a function (like Re, Im , or by default #&) which should be
        applied to the dipole list before the spectrum  is taken.";
ωPower::usage = "ωPower is an option for getSpectrum  which specifies a
        power of frequency which should multiply  the spectrum .";
DifferentiationOrder::usage = "DifferentiationOrderis an option for
        getSpectrum  which specifies the order to which the dipole
        list should be differentiatedbefore the spectrum  is taken.";
Protect[Polarization, ComplexPart , ωPower, DifferentiationOrder];

Begin["`Private`"];
Options[getSpectrum ] = {Polarization→ False, ComplexPart → (# &),
        ωPower → 0, DifferentiationOrder→ 0} ~Join~ standardOptions;

getSpectrum ::diffOrd = "Invalid differentiationorder `1`.";
getSpectrum ::ωPow = "Invalid ω power `1`.";

getSpectrum [dipoleList_, OptionsPattern[]] := Block[
    {polarizationVector, differentiatedList, depth, dimensions ,
      num  = OptionValue[TotalCycles],
      npp = OptionValue[PointsPerCycle], ω = OptionValue[CarrierFrequency], δt = 2π/ω/npp
    },
    polarizationVector= OptionValue[Polarization]/Norm [OptionValue[Polarization]];

    differentiatedList= OptionValue[ComplexPart ][Piecewise[{
          {dipoleList, OptionValue[DifferentiationOrder] == 0},
          {1/(2 δt) (Most[Most[dipoleList]] - Rest[Rest[dipoleList]]),
            OptionValue[DifferentiationOrder] == 1},
          {1/δt^2 (Most[Most[dipoleList]] - 2 Most[Rest[dipoleList]] + Rest[Rest[dipoleList]]),
            OptionValue[DifferentiationOrder] == 2}}],
        Message[getSpectrum ::diffOrd, OptionValue[DifferentiationOrder]];
        Abort[]
      ];

    If[NumberQ [OptionValue[ωPower]], Null;, Message[getSpectrum ::ωPow, OptionValue[ωPower]];
      Abort[] ];

    num Table[
      (ω/num k)^(2 OptionValue[ωPower]), {k, 1, Round[Length[differentiatedList]/2]}
    ] × If[
      OptionValue[Polarization] === False, (*unpolarized spectrum *)
      (*funky depth thing so this can take lists of numbers  and lists of vectors,
      of arbitrary length. Makes for easier benchmarking .*)
      depth = Length[Dimensions [dipoleList]];
```

```
        dimensions = If[Length[#] > 1, #[[2]], 1 (*#[[1]]*)] &[Dimensions [dipoleList]];
        Sum [Abs[
              Fourier[
                  If[depth > 1, Re[differentiatedList[[All, i]]], Re[differentiatedList[[All]]]]
                  , FourierParameters → {-1, 1}
                ][[1 ;; Round[Length[differentiatedList]/2]]]
            ]^2, {i, 1, dimensions }]
        , (*polarized spectrum *)
        Abs[
            Transpose[Table[
                    Fourier[
                      Re[differentiatedList[[All, i]]]
                      , FourierParameters → {-1, 1}
                    ]
                    , {i, 1, 2}]][[
                1 ;; Round[Length[differentiatedList/2]]].polarizationVector
          ]^2
      ]
  ]
End[];
```

spectrumPlotter

spectrumPlotter takes a spectrum and a list of options and returns a plot of the spectrum. The available options are

- a FrequencyAxis option, which will give the harmonic order as a horizontal axis by default, and an arbitrary scale with any other option,
- all the options of harmonicOrderAxis, which will be passed to the call that makes the horizontal axis, and
- all the options of ListLinePlot, which will be used to format the plot.

```
spectrumPlotter ::usage = "spectrumPlotter [spectrum ] plots
        the given spectrum  with an appropriate axis in a log10 scale.";
FrequencyAxis::usage = "FrequencyAxis is an option for spectrumPlotter
        which specifies the axis to use.";
Protect[FrequencyAxis];
Begin["`Private`"];
Options[spectrumPlotter ] = Join[{FrequencyAxis→ "HarmonicOrder "},
      Options[harmonicOrderAxis ], Options[ListLinePlot]];
spectrumPlotter [spectrum_ , options: OptionsPattern[]] := ListPlot[
    {Which[
          OptionValue[FrequencyAxis] === "HarmonicOrder ",
          harmonicOrderAxis ["TargetLength" → Length[spectrum ], Sequence@@FilterRules[
                {options} ~Join~ Options[spectrumPlotter ], Options[harmonicOrderAxis ]]],
          OptionValue[FrequencyAxis] === "Frequency",
          frequencyAxis["TargetLength" → Length[spectrum ], Sequence@@FilterRules[
                {options} ~Join~ Options[spectrumPlotter ], Options[harmonicOrderAxis ]]],
          True, Range[Length[spectrum ]]
        ],
        Log[10, spectrum ]
      }ᵀ
    , Sequence@@FilterRules[{options}, Options[ListLinePlot]]
    , Joined→ True
    , PlotRange → Full
    , PlotStyle → Thick
    , Frame  → True
    , Axes → False
    , ImageSize  → 800
  ]
End[];
```

biColorSpectrum

biColorSpectrum takes a time-dependent dipole list and produces overlaid plots of the right- and left-circular components of the spectrum, in red and blue respectively. It takes all the options of getSpectrum and spectrumPlotter, which are passed directly to the corresponding calls, as well as the options of Show, which can be used to modify the plot appearance.

```
Quit
```

```
biColorSpectrum ::usage =
    "biColorSpectrum [DipoleList] produces a two-colour spectrum  of DipoleList,
        separating the two circular polarizations.";
Begin["`Private`"];
Options[biColorSpectrum ] = Join[{PlotRange → All}, Options[Show],
      Options[spectrumPlotter ], DeleteCases[Options[getSpectrum ], Polarization→ False]];
biColorSpectrum [dipoleList_, options: OptionsPattern[]] := Show[{
      spectrumPlotter [
        getSpectrum [dipoleList, Polarization→ {1, +ⅈ},
          Sequence@@FilterRules[{options}, Options[getSpectrum ]]],
        PlotStyle → Red, Sequence@@FilterRules[{options}, Options[spectrumPlotter ]]],
      spectrumPlotter [
        getSpectrum [dipoleList, Polarization→ {1, -ⅈ},
          Sequence@@FilterRules[{options}, Options[getSpectrum ]]],
        PlotStyle → Blue, Sequence@@FilterRules[{options}, Options[spectrumPlotter ]]]
    }
    , PlotRange → OptionValue[PlotRange]
    , Sequence@@FilterRules[{options}, Options[Show]]
  ]
End[];
```

Various gate functions

Gate functions are used to suppress the contributions of extra-long trajectories with long excursion times, partly to reflect the effect of phase matching but mostly to keep integration times reasonable. They are provided to the main numerical integrator makeDipoleList via its Gate option.

```
SineSquaredGate::usage =
    "SineSquaredGate[nGateRamp ] specifies an integration gate with a sine-squared
        ramp , such that SineSquaredGate[nGateRamp ][ωt,nGate]
        has nGate flat periods and nGateRamp  ramp  periods.";
LinearRampGate ::usage = "LinearRampGate [nGateRamp ] specifies an integration gate
        with a linear ramp , such that SineSquaredGate[nGateRamp ][ωt,nGate]
        has nGate flat periods and nGateRamp  ramp  periods.";
Begin["`Private`"];
SineSquaredGate[nGateRamp_ ][ωτ_, nGate_] := Piecewise[{{1, ωτ ≤ 2 π (nGate - nGateRamp )},
      {Sin[(2 πnGate - ωτ)/(4 nGateRamp)]^2, 2 π (nGate - nGateRamp ) < ωτ ≤ 2 πnGate}, {0, nGate < ωτ}}]
LinearRampGate [nGateRamp_ ][ωτ_, nGate_] := Piecewise[{{1, ωτ ≤ 2 π (nGate - nGateRamp )},
      {-(ωτ - 2 π (nGate + nGateRamp ))/(2 πnGateRamp), 2 π (nGate - nGateRamp ) < ωτ ≤ 2 πnGate}, {0, nGate < ωτ}}]
End[];
```

getIonizationPotential

```
getIonizationPotential::usage =
    "getIonizationPotential[Target] returns the ionizationpotential
        of an atomic  target, e.g. \"Hydrogen\", in atomic  units.

getIonizationPotential[Target,q] returns the ionizationpotential
        of the q-th ion of the specified Target, in atomic  units.

getIonizationPotential[{Target,q}] returns the ionizationpotential
        of the q-th ion of the specified Target, in atomic  units.";
Begin["`Private`"];
getIonizationPotential[Target_, Charge_:0] :=
  UnitConvert[ElementData [Target, "IonizationEnergies"][[Charge+1]]/
      (Quantity[1, "AvogadroConstant"] Quantity[1, "Hartrees"])]
getIonizationPotential[{Target_, Charge_:0}] := getIonizationPotential[Target, Charge]
End[];
```

makeDipoleList: main numerical integrator

The main integration function is makeDipoleList, and its basic syntax is of the form makeDipoleList[VectorPotential→A]. Here the vector potential A must be a function object, such that for numeric t the construct A[t] returns a list of numbers after the appropriate field parameters have been introduced: thus the criterion is that, for a call of the form makeDipoleList[VectorPotential→A, FieldParameters→pars], a call of the form A[t]//.pars returns a list of numbers for numeric t. To see the available options use Options[makeDipoleList], and to get information on each option use the ?VectorPotential construct.

```
makeDipoleList::usage = "makeDipoleList [VectorPotential→A]
        calculates the dipole response to the vector potential A.";

VectorPotential::usage =
    "VectorPotential is an option for makeDipole  list which specifies the
        field's vector potential. Usage should be VectorPotential→A,
        where A[t]//.pars must  yield a list of numbers  for numeric
        t and parameters  indicated by FieldParameters →pars.";
VectorPotentialGradient::usage = "VectorPotentialGradient is an option for makeDipole
        list which specifies the gradient of the field's vector potential.
        Usage should be VectorPotentialGradient→GA, where GA[t]//.pars must
        yield a square matrix  of the same  dimension  as the vector potential
        for numeric  t and parameters  indicated by FieldParameters →pars.
        The indices must  be such that GA[t][[i,j]] returns ∂ᵢAⱼ[t].";
ElectricField::usage = "ElectricField is an option for makeDipole  list which specifies
        an electric field to use in the ionizationmatrix  element , in case the
        time  derivative of the vector potential is not desired. Usage should
        be ElectricField→F, where F[t]//.pars must  yield a list of numbers
        for numeric  t and parameters  indicated by FieldParameters →pars.";
FieldParameters::usage = "FieldParameters  is an option for makeDipole  list which ";
Preintegrals::usage =
    "Preintegrals is an option for makeDipole  list which specifies whether the preintegrals
        of the vector potential should be \"Analytic\" or \"Numeric \".";
ReportingFunction::usage = "ReportingFunction is an option for makeDipole  list
        which specifies a function used to report the results, either
        internally (by the default, Identity) or to an external file.";
Gate::usage = "Gate is an option for makeDipole  list which specifies the integration
        gate to use. Usage as Gate→g, nGate→n will gate the integral at
        time  ωt/ω by g[ωt,n]. The default is Gate→SineSquaredGate[1/2].";
nGate::usage = "nGate is an option for makeDipole  list which specifies
        the total number  of cycles in the integration gate.";
```

```
IonizationPotential::usage = "IonizationPotentialis an option for makeDipoleList
        which specifies the ionizationpotential Ip of the target.";
Target::usage = "Target is an option for makeDipoleList which specifies chemical
        species producing the HHG emission , pulling the ionization
        potential from the Wolfram ElementData curated data set.";
DipoleTransitionMatrixElement::usage = "DipoleTransitionMatrixElement is
        an option for makeDipoleList which secifies a function f to
        use as the dipole transition matrix element , or a pair of
        functions {fion, frec} to be used separately for the ionizationand
        recombination dipoels, to be used in the form f[p,κ]=f[p, √(2 Ip) ].";
ϵCorrection::usage = "ϵCorrection is an option for makeDipoleList which specifies the
        regularization correction ϵ, i.e. as used in the factor 1/(t-tt+iϵ)^(3/2).";
PointNumberCorrection::usage = "PointNumberCorrection is an option for makeDipoleList
        and timeAxis which specifies an extra number of points to be
        integrated over, which is useful to prevent Indeterminate errors
        when a Piecewise envelope is being differentiated at the boundaries.";
IntegrationPointsPerCycle::usage = "IntegrationPointsPerCycle is an option for
        makeDipoleList which controls the number of points per cycle to use for the
        integration. Set to Automatic , to follow PointsPerCycle, or to an integer.";
RunInParallel::usage = "RunInParallel is an option for makeDipoleList which controls
        whether each RB-SFA instance is parallelized. It accepts False as the
        (Automatic ) option, True, to parallelize each instance, or a pair of
        functions {TableCommand , SumCommand } to use for the iteration and
        summing , which could be e.g. {Inactive[ParallelTable], Inactive[Sum ]}.";
Simplifier::usage = "Simplifier is an option for makeDipoleList which specifies a
        function to use to simplify the intermediate and final analytical results.";
CheckNumericFields::usage = "CheckNumericFields is an option for
        makeDipoleList which specifies whether to check for
        numeric values of A[t] and GA[t] for numeric t.";
QuadraticActionTerms::usage = "QuadraticActionTerms is an option for makeDipoleList
        which specifies whether to use quadratic terms in ∇A^2 in the action.";

Protect[VectorPotential, VectorPotentialGradient, ElectricField, FieldParameters ,
    Preintegrals, ReportingFunction, Gate, nGate, IonizationPotential, Target, ϵCorrection,
    PointNumberCorrection , DipoleTransitionMatrixElement, IntegrationPointsPerCycle,
    RunInParallel, Simplifier, CheckNumericFields , QuadraticActionTerms ];


Begin["`Private`"];
Options[makeDipoleList ] = standardOptions~Join~{
        VectorPotential→Automatic , FieldParameters →{},
        VectorPotentialGradient→None, ElectricField→Automatic ,
        Preintegrals→"Analytic", ReportingFunction→Identity,
        Gate→SineSquaredGate[1/2], nGate→3/2, ϵCorrection→0.1,
        IonizationPotential→0.5,
        Target→Automatic , DipoleTransitionMatrixElement→hydrogenicDTME,
        PointNumberCorrection →0, Verbose→0, CheckNumericFields →True,
        RunInParallel →Automatic ,
        Simplifier→Identity, QuadraticActionTerms →True
      };
makeDipoleList::gate =
    "The integration gate g provided as Gate→`1` is incorrect. Its usage as
        g[`2`,`3`] returns `4` and should return a number .";
makeDipoleList::pot = "The vector potential A provided as VectorPotential→`1`
        is incorrect or is missing FieldParameters . Its usage as
        A[`2`] returns `3` and should return a list of numbers .";
makeDipoleList::efield = "The electric field f provided as ElectricField→`1` is incorrect
        or is missing FieldParameters . Its usage as F[`2`] returns `3` and should
```

```
        return a list of numbers . Alternatively, use ElectricField→Automatic .";
makeDipoleList::gradpot = "The vector potential GA provided as VectorPotentialGradient→`1`
        is incorrect or is missing  FieldParameters . Its usage as
        GA[`2`] returns `3` and should return a square matrix  of
        numbers . Alternatively, use VectorPotentialGradient→None.";
makeDipoleList::preint = "Wrong Preintegrals option `1`. Valid
        options are \"Analytic\" and \"Numeric \".";
makeDipoleList::runpar = "Wrong RunInParallel option `1`.";
makeDipoleList::carrfreq = "Non-numeric  option CarrierFrequency `1`.";
```

```
makeDipoleList[OptionsPattern[]] := Block[
    {
      num  = OptionValue[TotalCycles], npp = OptionValue[PointsPerCycle], ω,
      dipoleRec, dipoleIon, κ,
      A, F, GA, pi, ps, S,
      gate, tGate, setPreintegral,
      tInit, tFinal, δt, δtint, ϵ = OptionValue[ϵCorrection],
      AInt, A2Int, GAInt, GAdotAInt, AdotGAInt, GAIntInt,
      PScorrectionInt, constCorrectionInt, GAIntdotGAIntInt, QuadMatrix, q,
      simplifier, prefactor, integrand, dipoleList,
      TableCommand  , SumCommand
    },

    A[t_] = OptionValue[VectorPotential][t] //. OptionValue[FieldParameters ];
    If[
      OptionValue[ElectricField] === Automatic , F[t_] = -D[A[t], t];,
      F[t_] = OptionValue[ElectricField][t] //. OptionValue[FieldParameters ];
    ];
    GA[t_] = If[
        TrueQ[OptionValue[VectorPotentialGradient] == None],
        Table[0, {Length[A[tInit]]}, {Length[A[tInit]]}],
        OptionValue[VectorPotentialGradient][t] //. OptionValue[FieldParameters ]
      ];

    ω = OptionValue[CarrierFrequency];
    If[!NumberQ [ω] && TrueQ[OptionValue[CheckNumericFields ]],
      Message[makeDipoleList::carrfreq, ω];
      Abort[]];
    tInit= 0;
    tFinal= 2π/ω num ;
    (*looping timestep *)
    δt = (tFinal- tInit)/(num ×npp+ OptionValue[PointNumberCorrection ]);
    (*integration timestep *)
    δtint= If[OptionValue[IntegrationPointsPerCycle] === Automatic ,
        δt, (tFinal- tInit)/(num ×OptionValue[IntegrationPointsPerCycle] +
              OptionValue[PointNumberCorrection ])];
    tGate = OptionValue[nGate] 2π/ω;
    (*Check potential and potential gradient for correctness.*)
    (*To do: change logic conditions to constructions on VectorQ[#,NumberQ ]&
      and MatrixQ.*)
    If[TrueQ[OptionValue[CheckNumericFields ]],
```

```
  With[{ωtRandom = RandomReal [{ωtInit, ωtFinal}]},
    If[!And@@(NumberQ /@A[ωtRandom /ω]), Message[makeDipoleList::pot,
        OptionValue[VectorPotential], ωtRandom, A[ωtRandom]];
      Abort[]];
    If[!And@@(NumberQ /@Flatten[GA[ωtRandom /ω]]), Message[makeDipoleList::gradpot,
        OptionValue[VectorPotentialGradient], ωtRandom, GA[ωtRandom]];
      Abort[]];
    If[!And@@(NumberQ /@F[ωtRandom /ω]), Message[makeDipoleList::efield,
        OptionValue[ElectricField], ωtRandom, F[ωtRandom]];
      Abort[]];
  ];

gate[ωt_] := OptionValue[Gate][ωt, OptionValue[nGate]];
With[{ωtRandom = RandomReal [{ωtInit, ωtFinal}]},
  If[!TrueQ[NumberQ [gate[ωtRandom]]],
    Message[makeDipoleList::gate,
      OptionValue[Gate], ωtRandom, OptionValue[nGate], gate[ωtRandom]];
    Abort[]]
];

(*Target setup*)
Which[
  OptionValue[Target] === Automatic, κ = Sqrt[2 OptionValue[IonizationPotential]],
  True, κ = Sqrt[2 getIonizationPotential[OptionValue[Target]]]
];
With[{dim = Length[A[RandomReal [{ωtInit, ωtFinal}]]]},
  (*Explicit conjugation of the
    recombination matrix element to keep the integrand analytic.*)
  Which[
      Head[OptionValue[DipoleTransitionMatrixElement]] === List,
      dipoleIon[{p1_, p2_, p3_}[[1 ;; dim]], κκ_] =
        First[OptionValue[DipoleTransitionMatrixElement]][{p1, p2, p3}[[1 ;; dim]], κκ];
      dipoleRec[{p1_, p2_, p3_}[[1 ;; dim]], κκ_] =
        Assuming [{{p1, p2, p3, κκ} ∈ Reals}, Simplify [
            Conjugate[Last[OptionValue[DipoleTransitionMatrixElement]][
                {p1, p2, p3}[[1 ;; dim]], κκ]]
          ]];
      , True,
      dipoleIon[{p1_, p2_, p3_}[[1 ;; dim]], κκ_] =
        OptionValue[DipoleTransitionMatrixElement][{p1, p2, p3}[[1 ;; dim]], κκ];
      dipoleRec[{p1_, p2_, p3_}[[1 ;; dim]], κκ_] =
        Assuming [{{p1, p2, p3, κκ} ∈ Reals}, Simplify [
            Conjugate[OptionValue[DipoleTransitionMatrixElement][
                {p1, p2, p3}[[1 ;; dim]], κκ]]
          ]];
    ];
];


simplifier = OptionValue[Simplifier];
q = Boole[TrueQ[OptionValue[QuadraticActionTerms]]];

setPreintegral[integralVariable_, preintegrand_,
    dimensions_, integrateWithoutGradient_, parametric_] := Which[
```

```
    OptionValue[VectorPotentialGradient] =!= None || TrueQ[integrateWithoutGradient],
    (*Vector potential gradient specified,
    or integral variable does not depend on it, so integrate*)
    Which[
      OptionValue[Preintegrals] == "Analytic",
      integralVariable[t_, tt_] = simplifier[
            ((# /. {τ → t}) - (# /. {τ → tt})) &[Integrate[preintegrand[τ, tt], τ]]];

      , OptionValue[Preintegrals] == "Numeric ",
      Which[
        TrueQ[Not[parametric ]],
        Block[{innerVariable},
          integralVariable[t_, tt_] = (innerVariable[t] - innerVariable[tt] /. First[
                NDSolve[{innerVariable'[τ] == preintegrand[τ],
                  innerVariable[tInit] == ConstantArray[0, dimensions ]},
                 innerVariable, {τ, tInit, tFinal}, MaxStepSize → 0.25 / ω]
               ])
          ];
        , True,
        Block[{matrixpreintegrand , innerVariable, τpre},
          matrixpreintegrand [indices_, t_?NumericQ , tt_?NumericQ ] :=
            preintegrand[t, tt][[## & @@ indices]];
          integralVariable[t_, tt_] = Array[(
                  innerVariable[##][t - tt, tt] /. First@NDSolve[{
                    D[innerVariable[##][τpre, tt], τpre] ==
                      Piecewise[{{matrixpreintegrand [{##}, tt + τpre, tt],
                            tt + τpre ≤ tFinal}}, 0],
                    innerVariable[##][0, tt] == 0
                   }, innerVariable[##]
                   , {τpre, 0, tFinal - tInit}, {tt, tInit, tFinal}
                   , MaxStepSize → 0.25 / ω
                  ]
                ) &, dimensions ];
          ]
        ];
      ];
    , OptionValue[VectorPotentialGradient] === None,
    (*Vector potential gradient has not been specified,
    and integral variable depends on it, so return appropriate zero matrix *)
    integralVariable[t_] = ConstantArray[0, dimensions ];
    integralVariable[t_, tt_] = ConstantArray[0, dimensions ];
  ];

Apply[setPreintegral,

  ⎛AInt                A[#1] &
  ⎜A2Int               A[#1].A[#1] &
  ⎜GAInt               GA[#1] &
  ⎜GAdotAInt           GA[#1].A[#1] &
  ⎜AdotGAInt           A[#1].GA[#1] &
  ⎜GAIntInt            GAInt[#1, #2] &
  ⎜PScorrectionInt     GAdotAInt[#1, #2] + A[#1].GAInt[#1, #2] - q GAInt[#1, #2]^T.GAdotAInt[#1, #
  ⎜GAIntdotGAIntInt    q GAInt[#1, #2]^T.GAInt[#1, #2] &
  ⎝constCorrectionInt  (A[#1] - q/2 GAdotAInt[#1, #2]).GAdotAInt[#1, #2] &
```

```
    ),{1}];

(*{∫_t0^t A(τ)dτ, ∫_t0^t A(τ)^2 dτ, ∫_t0^t ∇A(τ)dτ, ∫_t0^t ∇A(τ)·A(τ)dτ, ∫_t0^t A(τ)·∇A(τ)dτ, ∫_t'^t ∫_t'^τ ∇A(τ')dτ'dτ,
    ∫_t'^t ∫_t'^τ ∂_j A_k(τ') A_k(τ')dτ' + A_k(τ) ∫_t'^τ ∂_k A_j(τ')dτ' - ∫_t'^τ ∂_i A_j(τ')dτ' ∫_t'^τ ∂_i A_k(τ') A_k(τ')dτ'dτ,
    ∫_t'^t ∫_t0^t ∂_i A_j(τ') A_j(τ')dτ' ∫_t0^t ∂_i A_k(τ') A_k(τ')dτ'dτ,
    ∫_t'^t (A_k(τ) - 1/2 ∫_t'^τ ∂_k A_i(τ') A_i(τ')dτ') · ∫_t'^τ ∂_k A_j(τ') A_j(τ')dτ'dτ};*)


(*Displaced momentum    *)
pi[p_, t_, tt_] := p + A[t] - GAInt[t, tt].p - GAdotAInt[t, tt];

(*Quadratic coefficient in nondipole action*)
QuadMatrix[t_, tt_] := (GAIntInt[t, tt] + GAIntInt[t, tt]^T)/2 - 1/2 GAIntdotGAIntInt[t, tt];

(*Stationary momentum    and action*)
ps[t_, tt_] := ps[t, tt] =
    -1/(t - tt - i ϵ) Inverse[IdentityMatrix[Length[A[tInit]]] - 1/(t - tt - i ϵ) 2 QuadMatrix[t, tt]] .
      (AInt[t, tt] - PScorrectionInt[t, tt]);
S[t_, tt_] := simplifier[
    1/2 (Total[ps[t, tt]^2] + κ^2) (t - tt) + ps[t, tt].AInt[t, tt] + 1/2 A2Int[t, tt] - (
      ps[t, tt].QuadMatrix[t, tt].ps[t, tt] +
        ps[t, tt].PScorrectionInt[t, tt] + constCorrectionInt[t, tt]
    )
  ];

prefactor[t_, τ_] := i (2π/(ϵ + i τ))^(3/2) dipoleRec[pi[ps[t, t - τ], t, t - τ], κ] ×
    dipoleIon[pi[ps[t, t - τ], t - τ, t - τ], κ].F[t - τ];
integrand[t_, τ_] := prefactor[t, τ] Exp[-i S[t, t - τ]] gate[ω τ];


(*Debugging constructs. Verbose→
  1 prints information about the internal functions. Verbose→
    2 returns all the relevant internal functions and stops. Verbose→
      3 for quantum -orbit constructs.*)
Which[
  OptionValue[Verbose] == 1, Information /@ {A, GA, ps, pi, S, AInt, A2Int, GAInt, GAdotAInt,
      AdotGAInt, GAIntInt, PScorrectionInt, constCorrectionInt, GAIntdotGAIntInt},
  OptionValue[Verbose] == 2, Return[With[{t = Symbol ["t"],
      tt = Symbol ["tt"], τ = Symbol ["τ"],
      p = {Symbol ["p1"], Symbol ["p2"], Symbol ["p3"]}[[1 ;; Length[A[ω tInit]]]]},
    {A[t], GA[t], ps[t, tt], pi[p, t, tt], S[t, tt], AInt[t, tt],
      A2Int[t, tt], GAInt[t, tt], GAdotAInt[t, tt], AdotGAInt[t, tt],
```

```
            GAIntInt[t, tt], PScorrectionInt[t, tt], constCorrectionInt[t, tt],
            GAIntdotGAIntInt[t, tt], QuadMatrix[t, tt], integrand[t, τ]}]],
      OptionValue[Verbose] == 3,
      Return[{
          Function[Evaluate[prefactor[#1, #1 - #2]]], Function[Evaluate[S[#1, #2]]]
        }]
    ];

    (*Single-run parallelization*)
    Which[
      OptionValue[RunInParallel] === Automatic ||
        OptionValue[RunInParallel] === False, TableCommand   = Table;
      SumCommand    = Sum ;,
      OptionValue[RunInParallel] === True, TableCommand   = ParallelTable;
      SumCommand    = Sum ;,
      True, TableCommand   = OptionValue[RunInParallel][[1]];
      SumCommand    = OptionValue[RunInParallel][[2]];
    ];

    (*Numerical  integration loop*)
    dipoleList= Table[
        OptionValue[ReportingFunction][
          δtintSum [(
                integrand[t, τ]
              ), {τ, 0,
                If[OptionValue[Preintegrals] == "Analytic", tGate, Min[t - tInit, tGate]], δtint}]
        ]
        , {t, tInit, tFinal, δt}
      ];
    dipoleList

  ]
End[];
```

Quantum orbit functions suite

Complex root finder

This section implements a routine for solving contains subroutines for the numerical solution of multiple simultaneous complex-valued transcendental equations, essentially by using the Newton's-method solver implemented in FindRoot, and seeding it multiple times with a random (or quasi-random) seed from a box. This code has been taken from the EPToolbox package, which is located and better documented at https://github.com/episanty/EPToolbox, and it is also documented in http://mathematica.stackexchange.com/a/57821/1000.

```
FindComplexRoots ::usage =
    "FindComplexRoots [e1==e2, {z, zmin , zmax }] attempts  to find complex  roots of the
        equation e1==e2 in the complex  rectangle with corners zmin  and zmax .

FindComplexRoots [{e1==e2, e3==e4, ...}, {z1, z1min , z1max },
        {z2, z2min , z2max }, ...] attempts  to find complex  roots of
        the given system  of equations in the multidimensional  complex
        rectangle with corners z1min , z1max , z2min , z2max , ....";
Seeds::usage = "Seeds is an option for FindComplexRoots  which determines  how many
        initialseeds are used to attempt  to find roots of the given equation.";
SeedGenerator::usage = "SeedGenerator is an option for FindComplexRoots  which determines
        the function  used to generate the seeds for the internal FindRoot
        call. Its value can be RandomComplex  , RandomNiederreiterComplexes  ,
        RandomSobolComplexes  , DeterministicComplexGrid , or any
        function f such that f[{zmin , zmax }, n] returns n complex
        numbers  in the rectancle with corners zmin  and zmax .";

Options[FindComplexRoots ] = Join[Options[FindRoot], {Seeds -> 50,
        SeedGenerator -> RandomComplex  , Tolerance -> Automatic , Verbose -> False}];
SyntaxInformation [FindComplexRoots ] = {"ArgumentsPattern " ->
        {_, {_, _, _}, OptionsPattern[]}, "LocalVariables" -> {"Table", {2, ∞}}};
FindComplexRoots ::seeds = "Value of option Seeds -> `1` is not a positive integer.";
FindComplexRoots ::tol =
    "Value of option Tolerance -> `1` is not Automatic  or a number  in [0,∞).";
$MessageGroups = Join[$MessageGroups, {"FindComplexRoots " :> {FindRoot::lstol}}]

Protect[Seeds];
Protect[SeedGenerator];

Begin["`Private`"];
FindComplexRoots [equations_List, domainSpecifiers__, ops : OptionsPattern[]] :=
  Block[{seeds, tolerances},
    If[! IntegerQ[Rationalize[OptionValue[Seeds]]] || OptionValue[Seeds] ≤ 0,
      Message[FindComplexRoots ::seeds, OptionValue[Seeds]]];
    If[! (OptionValue[Tolerance] === Automatic || OptionValue[Tolerance] ≥ 0),
      Message[FindComplexRoots ::tol, OptionValue[Seeds]]];

    seeds = OptionValue[SeedGenerator][{domainSpecifiers}[[All, {2, 3}]], OptionValue[Seeds]];
    tolerances = Which[
        ListQ[OptionValue[Tolerance]], OptionValue[Tolerance],
        True, ConstantArray[
          Which[
            NumberQ [OptionValue[Tolerance]], OptionValue[Tolerance],
            True, 10^If[NumberQ [OptionValue[WorkingPrecision]],
                2-OptionValue[WorkingPrecision], 2-$MachinePrecision]
          ]
          , Length[{domainSpecifiers}]]
      ];

    If[OptionValue[Verbose], Hold[], Hold[FindRoot::lstol]] /. {
        Hold[messageSequence___ ] :> Quiet[
            DeleteDuplicates[
              Select[
                Check[
                      FindRoot[
                        equations
                        , Evaluate[Sequence@@Table[{{domainSpecifiers}[[j, 1]],
                                #[[j]]}, {j, Length[{domainSpecifiers}]}]]]
```

```
                    , Evaluate[Sequence @@ FilterRules[{ops}, Options[FindRoot]]]
                  ],
                  ## &[]
                ] &/@seeds,
              Function[
                repList,
                ReplaceAll[
                  Evaluate[And@@Table[
                        And[
                          Re[{domainSpecifiers}[[j, 2]]] ≤ Re[
                               {domainSpecifiers}[[j, 1]]] ≤ Re[{domainSpecifiers}[[j, 3]]],
                          Im [{domainSpecifiers}[[j, 2]]] ≤ Im [{domainSpecifiers}[[
                                 j, 1]]] ≤ Im [{domainSpecifiers}[[j, 3]]]
                        ]
                        , {j, Length[{domainSpecifiers}]}]]
                  , repList]
                ]
              ],
              Function[{repList1, repList2},
                And@@Table[
                    Abs[({domainSpecifiers}[[j, 1]] /. repList1) -
                          ({domainSpecifiers}[[j, 1]] /. repList2)] < tolerances[[j]]
                    , {j, Length[{domainSpecifiers}]}]
              ]
            ]
            , {messageSequence }]}
  ]
FindComplexRoots [e1_ == e2_, {z_, zmin_ , zmax_ }, ops : OptionsPattern[]] :=
  FindComplexRoots [{e1 == e2}, {z, zmin , zmax }, ops]
End[];
```

Quasirandom number generators

This section implements quasirandom number generators for use with FindComplexRoots. As above, this code has been taken from the EPToolbox package, which is located and better documented at https://github.com/episanty/EP-Toolbox, and it is also documented in http://mathematica.stackexchange.com/a/57821/1000.

RandomSobolComplexes

```
RandomSobolComplexes  ::usage =
    "RandomSobolComplexes  [{zmin , zmax }, n] generates a low-discrepancy
        Sobol sequence of n quasirandom  complex  numbers
        in the rectangle with corners zmin  and zmax .

RandomSobolComplexes  [{{z1min ,z1max },{z2min ,z2max },...},n] generates a low-discrepancy
        Sobol sequence of n quasirandom  complex  numbers  in the multi -dimensional
        rectangle with corners {z1min ,z1max },{z2min ,z2max },....";
```

```
Begin["`Private`"];
RandomSobolComplexes [pairsList_, number_ ] := Map[
    Function[randomsList ,
      pairsList[[All, 1]] + Complex @@@Times [
            ReIm [pairsList[[All, 2]] - pairsList[[All, 1]]],
            randomsList
          ]
    ],
    BlockRandom [
      SeedRandom [Method → {"MKL", Method → {"Sobol", "Dimension " → 2 Length[pairsList]}}];
      SeedRandom [];
      RandomReal [{0, 1}, {number , Length[pairsList], 2}]
    ]
  ]
RandomSobolComplexes [{zmin_ ?NumericQ , zmax_ ?NumericQ }, number_ ] :=
  RandomSobolComplexes [{{zmin , zmax }}, number ][[All, 1]]
End[];
```

RandomNiederreiterComplexes

```
RandomNiederreiterComplexes ::usage =
    "RandomNiederreiterComplexes [{zmin , zmax }, n] generates
        a low-discrepancy Niederreiter sequence of n quasirandom
        complex numbers in the rectangle with corners zmin and zmax .

RandomNiederreiterComplexes [{{z1min ,z1max },{z2min ,z2max },...},n]
        generates a low-discrepancy Niederreiter sequence of n
        quasirandom complex numbers in the multi -dimensional
        rectangle with corners {z1min ,z1max },{z2min ,z2max },....";

Begin["`Private`"];
RandomNiederreiterComplexes [pairsList_, number_ ] := Map[
    Function[randomsList ,
      pairsList[[All, 1]] + Complex @@@Times [
            ReIm [pairsList[[All, 2]] - pairsList[[All, 1]]],
            randomsList
          ]
    ],
    BlockRandom [
      SeedRandom [
        Method → {"MKL", Method → {"Niederreiter", "Dimension " → 2 Length[pairsList]}}];
      SeedRandom [];
      RandomReal [{0, 1}, {number , Length[pairsList], 2}]
    ]
  ]
RandomNiederreiterComplexes [{zmin_ ?NumericQ , zmax_ ?NumericQ }, number_ ] :=
  RandomNiederreiterComplexes [{{zmin , zmax }}, number ][[All, 1]]
End[];
```

DeterministicComplexGrid

```
DeterministicComplexGrid ::usage =
    "DeterministicComplexGrid [{zmin , zmax }, n] generates a grid of about n equally
        spaced complex  numbers  in the rectangle with corners zmin  and zmax .

DeterministicComplexGrid [{{z1min ,z1max },{z2min ,z2max },...},n] generates a regular grid
        of about n equally spaced complex  numbers  in the multi -dimensional
        rectangle with corners {z1min ,z1max },{z2min ,z2max },....";

Begin["`Private`"];
DeterministicComplexGrid [pairsList_, number_ ] :=
  Block[{sep, separationsList, gridPointBasis, k},
    sep = NestWhile[0.99 # &, Min[Flatten[ReIm [pairsList[[All, 2]] - pairsList[[All, 1]]]]],
        Times @@ 1/(0.99 #) Floor[Flatten[ReIm [pairsList[[All, 2]] - pairsList[[All, 1]]]], 0.99 #] ≤
          number &];
    separationsList = Round[1/sep Floor[Flatten[ReIm [pairsList[[All, 2]] - pairsList[[All, 1]]]],
      sep]];
    gridPointBasis = MapThread[
        Function[{l, n}, Range[l[[1]], l[[2]], (l[[2]] - l[[1]])/(n + 1)][[2 ;; -2]]],
        {Flatten[Transpose[ReIm [pairsList], {1, 3, 2}], 1], separationsList}
      ];
    Flatten[Table[
        Table[k[2 j - 1] + i k[2 j], {j, 1, Length[pairsList]}],
        Evaluate[Sequence @@ Table[{k[j], gridPointBasis[[j]]}, {j, 1, 2 Length[pairsList]}]]
      ], Evaluate[Range[1, 2 Length[pairsList]]]]
  ]
DeterministicComplexGrid [{zmin_ ?NumericQ , zmax_ ?NumericQ }, number_ ] :=
  DeterministicComplexGrid [{{zmin , zmax }}, number ][[All, 1]]
End[];
```

RandomComplex

Updating RandomComplex to handle input of the form RandomComplex[{{0, 1+*i*}, {2, 3+*i*}}, n].

```
Begin["`Private`"];
Unprotect[RandomComplex ];
RandomComplex [{range1_List, moreRanges___ }, number_ ] :=
  Transpose[RandomComplex [#, number ] & /@ {range1, moreRanges }]
Protect[RandomComplex ];
End[];
```

The following code places this redefinition as an initialization code for any parallelized subkernels that may get launched later (cf. mm.se/q/131856). This version, in addition, checks whether there is already any code in $InitCode and, if there is, it appends its own code there.

```
Parallelize;
If[Head[Parallel`Developer`$InitCode] =!= Hold,
    Parallel`Developer`$InitCode= Hold[]
  ];
Parallel`Developer`$InitCode= Join[
      Parallel`Developer`$InitCode,
      Hold[
        Unprotect[RandomComplex  ];
        RandomComplex  [
            {Private`range1_List, Private`moreRanges___ }, Private`number_ ] := Transpose[
            RandomComplex  [#, Private`number ] &/@{Private`range1, Private`moreRanges }];
        Protect[RandomComplex  ];
      ]
    ];
```

GetSaddlePoints

```
GetSaddlePoints::usage =
    "GetSaddlePoints[Ω,S,{tmin ,tmax },{τmin ,τmax }] finds a list of solutions {t,τ}
        of the HHG temporal  saddle-point equations at harmonic  energy Ω
        for action S, in the range {tmin , tmax } of recombination time  and
        {τmin , τmax } of excursion time , where both ranges should be the
        lower-left and upper-right corners of rectangles in the complex  plane.

GetSaddlePoints[ΩRange,S,{tmin ,tmax },{τmin ,τmax }] finds solutions of
        the HHG temporal  saddle-point equations for a range of harmonic
        energies ΩRange, and returns an Association with each harmonic
        energy Ω indexing a list of saddle-point solution pairs {t,τ}.

GetSaddlePoints[Ωspec,S,{{{tmin 1,tmax 1},{τmin 1,τmax 1}},{{tmin 2,tmax 2},{τmin 2,τmax 2}},
        ...}] uses multiple  time  domains  and combines  the solutions.

GetSaddlePoints[Ωspec,S,{{urange,vrange},...},IndependentVariables→{u,v}] uses the explicit
        independent variables u and v to solve the equations and over the given ranges,
        where u and v can be any of \"RecombinationTime \", \"IonizationTime\"
        and \"ExcursionTime \", or their shorthands \"t\", \"tt\" and \"τ\" resp.";
SortingFunction::usage = "SortingFunction is an option of GetSaddlePoints
        which sets a function f, to be used as f[t,τ,S,Ω], to be
        used to sort the solutions, or a list of such functions.";
SelectionFunction::usage = "SelectionFunction is an option of GetSaddlePoints
        that sets a function f, to be used as f[t,τ,S,Ω],
        such that roots are only kept if f returns True.";
IndependentVariables::usage = "IndependentVariables is an option for GetSaddlePoints
        that specifies the two independent variables, out of \"RecombinationTime \",
        \"IonizationTime\" and \"ExcursionTime \" (or their shorthands
        \"t\", \"tt\" and \"τ\", respectively), to be used in solving the
        saddle-point equations, and which range over the given regions.";

FiniteDifference::usage =
    "FiniteDifferenceis a value for the option Jacobian of FindRoot, FindComplexRoots ,
        GetSaddlePoints, and related functions, which specifies that the Jacobian at
        each step should be evaluated using numerical  finite difference procedures.";

GetSaddlePoints::error = "Errors encountered for harmonic  energy Ω=`1`.";

Begin["`Private`"];
Options[GetSaddlePoints] =
    Join[{SortingFunction→ (#2 &), SelectionFunction→ (True &), IndependentVariables→
          {"RecombinationTime ", "ExcursionTime "}}, Options[FindComplexRoots ]];
```

```
Protect[SortingFunction, SelectionFunction, IndependentVariables, FiniteDifference];

GetSaddlePoints[Ωspec_, S_, {tmin_ , tmax_ }, {τmin_ , τmax_ }, options:OptionsPattern[]] :=
  GetSaddlePoints[Ωspec, S, {{{tmin , tmax }, {τmin , τmax }}}, options]

GetSaddlePoints[Ω_, S_, timeRanges_ , options:OptionsPattern[]] :=
  Block[{equations, roots, t = Symbol ["t"], tt = Symbol ["tt"],
      τ = Symbol ["τ"], indVars, depVar, depVarRule, tolerances},
    indVars = OptionValue[IndependentVariables] /.
        {"RecombinationTime " → "t", "ExcursionTime " → "τ", "IonizationTime" → "tt"};
    depVar = First[DeleteCases[{"t", "τ", "tt"}, Alternatives@@indVars]];
    depVarRule = depVar /. {"tt" → {tt → t - τ}, "t" → {t → tt + τ}, "τ" → {τ → t - tt}};
    equations = {D[S[t, tt], t] == Ω, D[S[t, tt], tt] == 0} /. depVarRule;
    tolerances = Which[
        ListQ[OptionValue[Tolerance]], OptionValue[Tolerance],
        True, ConstantArray[
          Which[
            NumberQ [OptionValue[Tolerance]], OptionValue[Tolerance],
            True, 10^If[NumberQ [OptionValue[WorkingPrecision]],
                2 - OptionValue[WorkingPrecision], 2 - $MachinePrecision]
          ]
          , 2]];

    SortBy[
      DeleteDuplicates[
        Flatten[Table[
            Select[
              Check[
                roots = ({t, τ} /. depVarRule) /. (FindComplexRoots [
                          equations
                          , Evaluate[Sequence[{Symbol [indVars[[1]]], range[[1, 1]], range[[1,
                                  2]]}, {Symbol [indVars[[2]]], range[[2, 1]], range[[2, 2]]}]]
                          , Evaluate[Sequence@@FilterRules[{options},
                                Options[FindComplexRoots ]]]
                          , SeedGenerator → RandomSobolComplexes
                          , Seeds → 50
                        ] /. {{} → (({t, τ} /. depVarRule) → {})})
                      (*to deal with empty  results*)
                , Message[GetSaddlePoints::error, Ω]; roots
              ]
              , Function[timesPair ,
                OptionValue[SelectionFunction][timesPair [[1]], timesPair [[2]], S, Ω]]
            ]
            , {range, timeRanges }], 1]
        , Function[{timesPair1 , timesPair2 },
          And@@Thread[Abs[timesPair1 - timesPair2 ] < tolerances] ]
      ]
      , If[
        ListQ[OptionValue[SortingFunction]],
        Table[Function[timesPair , f[timesPair [[1]], timesPair [[2]], S, Ω]],
          {f, OptionValue[SortingFunction]}],
        Function[timesPair , OptionValue[SortingFunction][timesPair [[1]], timesPair [[2]], S, Ω]]
      ]
    ]
  ]
GetSaddlePoints[ΩRange_List, S_, timeRanges_ , options:OptionsPattern[]] :=
```

```
  Association[ParallelTable[
      Ω→GetSaddlePoints[Ω, S, timeRanges , options]
      , {Ω, Sort[ΩRange]}]]

End[];
```

GetSaddlesFromSeeds

```
GetSaddlesFromSeeds ::usage =
    "GetSaddlesFromSeeds [{{t₁,τ₁},{t₂,τ₂},...},Ω,S] finds a list of solutions {t,τ}
        of the HHG temporal  saddle-point equations at harmonic  energy
        Ω for action S, using the given {tᵢ,τᵢ} as seeds for the process.

GetSaddlesFromSeeds [<|Ω₁→{{t₁₁,τ₁₁},{t₁₂,τ₁₂},...},Ω₂→{{t₂₁,τ₂₁},{t₂₂,τ₂₂},...},...|>,Ω,S]
        finds solutions of the HHG temporal  saddle-point equations,
        using the seeds list from  the Ωᵢ that's closest to
        Ω, or as specified by the value of KeyChooserFunction.

GetSaddlesFromSeeds [seeds,{Ω₁,Ω₂,...},S]
        iterates over the given set of harmonic  energies.";

SeedsChooserFunction::usage =
    "SeedsChooserFunction is an option for GetSaddlesFromSeeds  that
        specifies a function f (set by default to Nearest) that, when
        used as f[{Ω₁,Ω₂,...},Ω], should return the indices {Ωᵢ,Ωⱼ,...}
        corresponding to the seed sets {{{t_i1,τ_i1},...},{{t_j1,τ_j1},...}}
        to be used to solve the HHG saddle-point equations.";
RecalculateRoots::usage = "RecalculateRoots is an option for GetSaddlesFromSeeds
        that specifies whether to re-solve the saddle-point equations if
        the given harmonic  energy Ω is among  the set of keys of the given
        seeds association  The default is False, which is appropriate for S
        being the same  action used to find the seeds, in which case setting
        RecalculateRoots→True will produce multiple  FindRoot errors.  If
        using a different action than used to find the seeds, set to True.";

GetSaddlesFromSeeds ::error = "Errors encountered for harmonic  energy Ω=`1`.";
GetSaddlesFromSeeds ::norecalc =
    "Skipping re-calculation of roots at harmonic  energy `1` since it is
        already in the key set of the given seeds association  To run
        the calculation for this case set RecalculateRoots to True.";

Begin["`Private`"];
Options[GetSaddlesFromSeeds ] =
    Join[{RecalculateRoots→False, SeedsChooserFunction→Nearest}, Options[GetSaddlePoints]];
Protect[SeedsChooserFunction, RecalculateRoots];


GetSaddlesFromSeeds [seedsSpec_, ΩRange_List, S_, options:OptionsPattern[]] :=
  Association[ParallelTable[
      Ω→GetSaddlesFromSeeds [seedsSpec, Ω, S, options]
      , {Ω, Sort[ΩRange]}]]

GetSaddlesFromSeeds [seedsAssociation_Association,Ω_, S_, options:OptionsPattern[]] :=
  With[{keys = OptionValue[SeedsChooserFunction][Keys[seedsAssociation], Ω]},
    If[MemberQ [keys, Ω] && TrueQ[!OptionValue[RecalculateRoots]],
      Message[GetSaddlesFromSeeds ::norecalc, Ω];
      Return[seedsAssociation[Ω]]];
    GetSaddlesFromSeeds [Flatten[Values[seedsAssociation[Key/@keys]], 1], Ω, S, options]
  ]
```

```
GetSaddlesFromSeeds [seedsList_List, Ω_?NumberQ , S_, options:OptionsPattern[]] := Block[
    {equations, roots, t = Symbol ["t"], tt = Symbol ["tt"],
      τ = Symbol ["τ"], indVars, depVar, depVarRule, fullSeedVars, tolerances},
    indVars= OptionValue[IndependentVariables] /.
        {"RecombinationTime " → "t", "ExcursionTime " → "τ", "IonizationTime" → "tt"};
    depVar = First[DeleteCases[{"t", "τ", "tt"}, Alternatives@@indVars]];
    depVarRule = depVar /. {"tt" → {tt → t - τ}, "t" → {t → tt + τ}, "τ" → {τ → t - tt}};
    fullSeedVars[seed_] := ⟨|"t" → seed[[1]], "τ" → seed[[2]], "tt" → seed[[1]] - seed[[2]]|⟩;
    equations= {D[S[t, tt], t] == Ω, D[S[t, tt], tt] == 0} /. depVarRule;
    tolerances = Which[
        ListQ[OptionValue[Tolerance]], OptionValue[Tolerance],
        True, ConstantArray[
          Which[
            NumberQ [OptionValue[Tolerance]], OptionValue[Tolerance],
            True, 10^If[NumberQ [OptionValue[WorkingPrecision]],
                2 - OptionValue[WorkingPrecision], 2 - $MachinePrecision]
          ]
          , 2]];

    SortBy[
      DeleteDuplicates[
        Select[
          Table[
            Check[
              roots = ({t, τ} /. depVarRule) /. (
                    FindRoot[
                        equations
                        , {Symbol [#], fullSeedVars[seed][[#]]} & /@indVars
                        , Evaluate[Sequence@@FilterRules[{options}, Options[FindRoot]]]
                      ]
                      /. {{} → (({t, τ} /. depVarRule) → {})})
              , Message[GetSaddlesFromSeeds ::error, Ω]; roots
            ]
            , {seed, seedsList}]
          , Function[timesPair ,
            OptionValue[SelectionFunction][timesPair [[1]], timesPair [[2]], S, Ω]]
        ]
        , Function[{timesPair1 , timesPair2 },
          And@@Thread[Abs[timesPair1 - timesPair2 ] < tolerances] ]
      ]
      , If[
        ListQ[OptionValue[SortingFunction]],
        Table[Function[timesPair , f[timesPair [[1]], timesPair [[2]], S, Ω]],
          {f, OptionValue[SortingFunction]}],
        Function[timesPair , OptionValue[SortingFunction][timesPair [[1]], timesPair [[2]], S, Ω]]
      ]
    ]
  ]
End[];
```

ClassifyQuantumOrbits

```
ClassifyQuantumOrbits::usage =
    "ClassifyQuantumOrbits[saddlePoints,f] sorts an indexed set of saddle points
        of the form  <|Ω1→{{t11,τ11},{t12,τ12},...}...|> using a function f,
        which should turn f[t,τ,Ω] into an appropriate label, and returns an
        association of the form  <|label1→<|Ω1→<|1→{t,τ},2→{t,τ},...|>,...|>,...|>.

ClassifyQuantumOrbits[saddlePoints,f,sortFunction] uses
        the function sortFunction to sort the sets of saddle points
        {{t11,τ11},{t12,τ12},...} for each label and harmonic  energy.

ClassifyQuantumOrbits[saddlePoints,f,sortFunction,DiscardedLabels→{label1,label2,...}]
        specifies a list of labels to discard from  the final output.";
DiscardedLabels::usage = "DiscardedLabels is an option for ClassifyQuantumOrbits
        which specifies a list of labels to discard from  the final output.";

Begin["`Private`"];

Options[ClassifyQuantumOrbits] = {DiscardedLabels→{}};
Protect[DiscardedLabels];

ClassifyQuantumOrbits[saddlePointList_,
    classifierFunction_,sortingFunction_:Sort,OptionsPattern[]] := Map[
    Composition[
      Association,
      MapIndexed[#2[[1]]→#1&],
      sortingFunction
    ],
    Delete[
      AssociationTranspose[
        MapIndexed[
            GroupBy[classifierFunction@@#&][
                Flatten/@Transpose[{#1,ConstantArray[#2[[{1},1]],Length[#1]]}]]&
            ,saddlePointList][[All,All,All,{1,2}]]
      ]
      ,List/@OptionValue[DiscardedLabels]]
    ,{2}]

End[];
```

ReperiodSaddles

```
ClearAll[ReperiodSaddles]
ReperiodSaddles::usage =
    "ReperiodSaddles[{{t1,τ1},{t2,τ2},...},f] readjusts the assigned cycle of
        the saddle points {ti,τi}, returning the list {{t1+f[t1,τ1],τ1},...}.

ReperiodSaddles[<|Ω1→{{t11,τ11},...},Ω2→...|>,f] reperiods
        saddle-point pairs in a harmonic -energy-indexed association.

ReperiodSaddles[<|label1→<|Ω1→{{t11,τ11},...},...|>,...|>,f] reperiods
        saddle-point pairs of a classified set of saddle points.";

Begin["`Private`"];

ReperiodSaddles[pair_/;Depth[pair]==2,f_] := {pair[[1]]+f[pair[[1]],pair[[2]]],pair[[2]]}
ReperiodSaddles[association_,f_] := Apply[f,association,{Depth[association]-2}]

End[];
```

HessianRoot

```
HessianRoot::usage = "HessianRoot[S,t,τ] calculates the Hessian root √((2π)^2/(i^2 Det[∂^2_{t,tt} S])).";

Begin["`Private`"];

HessianRoot[S_, t_, τ_] := √(2π/(i Derivative[0, 2][S][t, t-τ])) √((2π Derivative[0, 2][S][t, t-τ])/(i (Derivative[2, 0][S][t, t-τ] Derivative[0, 2][S][t, t-τ] - Derivative[1, 1][S][t, t-τ]^2)))
End[];
```

FindStokesTransitions

```
FindStokesTransitions::usage =
    "FindStokesTransitions[S,<|Ω1→<|1→{t11,τ11},2→{t12,τ12}|>,Ω2→<|1→{t21,τ21},2→{t22,τ22}|>,...|>
    ] finds the set {{ΩS},{ΩAS},n} of the Stokes and
    anti-Stokes transition energies for the given set of saddle points,
    where Re(S) changes sign after the ΩS and Im (S) changes sign after
    the ΩAS, and n is the index of the member of the pair that should
    be chosen after the transition (taken as the member with a positive
    imaginary part of the action at the largest Ωi in the given keys).

FindStokesTransitions[S,<|label1→<|Ω1→...|>|>] finds the Stokes transitions for the given set
    of saddle-point curve pairs, and returns them labeled with the labeli.";
FindStokesTransitions::saddleno = "FindStokesTransitionscalled with `1` of `2`
    saddle-point sets of length different from 2, with set length
    structure `3`. Excluding those sets from the calculation.";
FindStokesTransitions::multipleS = "FindStokesTransitionsfound multiple
    Stokes transitions; using `1` to return a single transition.";
FindStokesTransitions::multipleAS = "FindStokesTransitionsfound multiple
    anti-Stokes transitions; using `1` to return a single transition.";
ChooserFunction::usage = "ChooserFunction is an option for FindStokesTransitionsthat
    specifies which transition to take if there are multiple transitions in
    the given dataset. The default is Last and gives the one with higher
    energy; to get the full set of transitions found use Full or Identity.";
ReperiodingFunction::usage = "ReperiodingFunctionis an option for FindStokesTransitions,
    SPAdipole and UAdipole which specifies a function f[t,τ] of recombination
    time t and excursion time τ that will be used to re-period the
    pairs {t,τ} into the form {t+f[t,τ],τ}. The default is Function[0],
    but if pairs are split it can be useful to set ReperiodingFunctionto
    Function[{t,τ},Floor[-Re[t-τ],2π/ω]] for ω the carrier frequency. In general,
    however, it is preferable to do this in a single go using ReperiodSaddles.";

Begin["`Private`"];

Protect[ReperiodingFunction, ChooserFunction];
Options[FindStokesTransitions] =
    {ReperiodingFunction→ Function[{t, τ}, 0], ChooserFunction→ Automatic };

FindStokesTransitions[S_,
    deeperAssociation_/; Depth[deeperAssociation] == 5, options:OptionsPattern[]] := Map[
    FindStokesTransitions[S, #, options] &,
    deeperAssociation
  ]
```

```
FindStokesTransitions[S_, saddlesAssociation_, options: OptionsPattern[]] :=
  Block[{reducedSaddlesAssociation, actionList, signsList, s, processor},
    reducedSaddlesAssociation = KeySort[Select[saddlesAssociation, Length[#] == 2 &]];
    If[Length[saddlesAssociation] - Length[reducedSaddlesAssociation] > 0,
      Message[FindStokesTransitions::saddleno,
        Length[saddlesAssociation] - Length[reducedSaddlesAssociation],
        Length[saddlesAssociation], First/@Tally/@Split[Values[Length/@saddlesAssociation]]
      ]
    ];
    actionList = ReIm [
        Map[(*reduces each Ω→<|1→S1,2→S2|> to Ω→(S1-S2)*)
          Apply[Subtract],
          MapIndexed[(*reduces each Ω→<|1→{t1,τ1},2→{t2,τ2}|> to Ω→<|1→S1,2→S2|>*)
            With[{t = #1[[1]] + OptionValue[ReperiodingFunction][#1[[1]], #2[[2]]],
                  τ = #1[[2]], Ω = #2[[1, 1]]},
              S[t, t - τ] - Ωt
            ] &
            , reducedSaddlesAssociation, {2}
          ]]
      ];
    signsList = Sign[Times [
          Rest[actionList],
          AssociationThread[Rest[Keys[actionList]], Most[Values[actionList]]]
        ]];
    processor = OptionValue[ChooserFunction] /. {Automatic → Last, Full → Identity};
    If[Length[Keys[Select[signsList, #[[1]] < 0 &]]] > 1,
      Message[FindStokesTransitions::multipleS, processor]];
    If[Length[Keys[Select[signsList, #[[2]] < 0 &]]] > 1,
      Message[FindStokesTransitions::multipleAS, processor]];
    {
      processor[Keys[Select[signsList, #[[1]] < 0 &]] /. {{} → {Missing["No transition"]}}],
      processor[Keys[Select[signsList, #[[2]] < 0 &]] /. {{} → {Missing["No transition"]}}],
      Sign[Last[actionList][[2]] /. {1 → 2, -1 → 1}
    }
  ]

End[];
```

SPAdipole

```
SPAdipole::usage =
    "SPAdipole[S,prefactor,Ω,{t,τ}] returns the saddle-point approximation amplitude
        corresponding to action S[t,t-τ]-Ωt and the given prefactor[t,t-τ].

SPAdipole[S,prefactor,Ω,<|1→{t1,τ1},2→{t2,τ2},...|>]
        returns the total harmonic -dipole contribution in the
        saddle-point approximation from the specified saddle points.

SPAdipole[S,prefactor,Ω,<|1→{t1,τ1},2→{t2,τ2}|>,transition] uses the given Stokes transition
        set to drop the relevant saddle after the anti-Stokes transition.";
SPAdipole::wrongno = "SPAdipole called with a Stokes transition but with
        an input association of length `1` at harmonic
        energy Ω=`2`. Reverting to unstructured evaluation.";
SPAdipole::invldtrns = "SPAdipole called with invalid Stokes transition
        set `1`. Reverting to unstructured evaluation.";

Begin["`Private`"];

Options[SPAdipole] = {ReperiodingFunction→ Function[{t, τ}, 0]};

SPAdipole[S_, prefactor_, Ω_, {t_, τ_}, options:OptionsPattern[]] :=
  Block[{tr = t + OptionValue[ReperiodingFunction][t, τ]},
    HessianRoot[S, tr, τ] prefactor[tr, tr - τ] Exp[-i S[tr, tr - τ] + i Ω tr]
  ]
SPAdipole[S_, prefactor_, Ω_, times_Association, options:OptionsPattern[]] := Block[{},
    Total[SPAdipole[S, prefactor, Ω, #, options] & /@ times ]
  ]

SPAdipole[S_, prefactor_, Ω_, times_Association,
    transition_, options:OptionsPattern[]] := Block[{},
    If[!NumberQ [transition[[2]]], Message[SPAdipole::invldtrns, transition];
      Return[SPAdipole[S, prefactor, Ω, times ]]];
    If[Length[times ] ≠ 2, Message[SPAdipole::wrongno, Length[times ], Ω];
      Return[SPAdipole[S, prefactor, Ω, times , options]]];
    If[Ω < transition[[2]],
      SPAdipole[S, prefactor, Ω, times , options],
      SPAdipole[S, prefactor, Ω, KeySelect[times , # == transition[[3]] &], options]
    ]
  ]

End[];
```

UAdipole

```
UAdipole::usage =
    "UAdipole[S,prefactor,Ω,<|1→{t1,τ1},2→{t2,τ2},...|>,transition] returns the total
        harmonic -dipole contribution in the uniform approximation from the
        specified saddle points, using the action S[t,t-τ]-Ωt and prefactor[t,t-τ],
        and taking the given Stokes transition set as a reference.";
UAdipole::saddleno = "UAdipole called with `1` time pairs at Ω=`2`.
        Reverting to the saddle-point approximation for this set.";
UAdipole::invldtrns = "UAdipole called with invalid Stokes transition set `1`.
        Reverting to the saddle-point approximation for this set.";

Begin["`Private`"];

Options[UAdipole] = {ReperiodingFunction→ Function[{t, τ}, 0]};

UAdipole[S_, prefactor_, Ω_, times_ , transition_, options:OptionsPattern[]] := (
    If[Length[times ] ≠ 2, Message[UAdipole::saddleno, Length[times ], Ω];
      Return[SPAdipole[S, prefactor, Ω, times ]]];
    If[!NumberQ [transition[[2]]], Message[UAdipole::invldtrns, transition];
      Return[SPAdipole[S, prefactor, Ω, times ]]];
    Block[
      {A1, A2, S1, S2, Ss, Sm , z,
        t1 = times [1][[1]] + OptionValue[ReperiodingFunction][times [1][[1]], times [1][[2]]],
        τ1 = times [1][[2]],
        t2 = times [2][[1]] + OptionValue[ReperiodingFunction][times [2][[1]], times [2][[2]]],
        τ2 = times [2][[2]]},
      A1 = HessianRoot[S, t1, τ1] prefactor[t1, t1 - τ1];
      S1 = S[t1, t1 - τ1] - Ω t1;
      A2 = HessianRoot[S, t2, τ2] prefactor[t2, t2 - τ2];
      S2 = S[t2, t2 - τ2] - Ω t2;
      Ss = (S1 + S2)/2; Sm = (S1 - S2)/2;
      If[Ω < transition[[2]], z = (-3/2 Sm)^(2/3),
        z = (-3/2 Sm)^(2/3) Exp[i (transition[[3]] /. {2 → -1, 1 → 1}) 2π/3]];
      Sqrt[6 π Sm] Exp[-i Ss + i π/4] ((A1 - i A2)/2 AiryAi[-z]/Sqrt[z] + i (A1 + i A2)/2 AiryAi'[-z]/z)
    ])

End[];
```

Package closure

End of package

```
EndPackage[];
```

Add to distributed contexts.

```
DistributeDefinitions["RBSFA`*"];
```