

LPAIR++  
0.1

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## 1 Todo List

Global `GamGam::GamGam` (`const unsigned int ndim_`, `int nOpt_`, `double x_[]`)

Figure out how this `nOpt_` parameter is affecting the final cross-section computation and events generation

## 2 Hierarchical Index

### 2.1 Class Hierarchy

This inheritance list is sorted roughly, but not completely, alphabetically:

<b>Event</b>	<b>3</b>
<b>GamGam</b>	<b>4</b>
<b>GamGamKinematics</b>	<b>7</b>
<b>InputParameters</b>	<b>11</b>
<b>MCGen</b>	<b>15</b>
<b>Particle</b>	<b>16</b>
<b>InelasticParticle</b>	<b>8</b>
<b>Vegas</b>	<b>19</b>

## 3 Data Structure Index

### 3.1 Data Structures

Here are the data structures with brief descriptions:

<b>Event</b>	
Kinematic information on the particles in the event	<b>3</b>
<b>GamGam</b>	
Computes the matrix element for a $\gamma\gamma \rightarrow \ell^+\ell^-$ process	<b>4</b>
<b>GamGamKinematics</b>	
List of kinematic cuts to apply on the central and outgoing phase space	<b>7</b>
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Kinematics of one particle	<b>16</b>
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Vegas Monte-Carlo integrator instance	<b>19</b>

## 4 Data Structure Documentation

### 4.1 Event Class Reference

Kinematic information on the particles in the event.

#### Public Member Functions

- void [Dump](#) ()
- [Particle](#) \* [GetByRole](#) (int *role\_*)
- int [SetParticle](#) ([Particle](#) \**part\_*)
- void [Store](#) (std::ofstream \*, double *weight\_*=1.)
- void [StoreLHERRecord](#) (std::ofstream \**of\_*, const double *weight\_*=1.)

*Stores the LHE block for this event.*

#### 4.1.1 Detailed Description

Class containing all the information on the in- and outgoing particles' kinematics

#### 4.1.2 Member Function Documentation

##### 4.1.2.1 void Event::Dump ( )

Dumps all the known information on every [Particle](#) object contained in this [Event](#) container in the output stream

##### 4.1.2.2 Particle\* Event::GetByRole ( int *role\_* )

Returns the pointer to the [Particle](#) object corresponding to a certain role in the process kinematics

#### Parameters

<i>role_</i>	The role the particle has to play in the process
--------------	--

#### Returns

A pointer to the requested [Particle](#) object

##### 4.1.2.3 int Event::SetParticle ( [Particle](#) \* *part\_* )

Sets the information on one particle in the process

#### Parameters

<i>part_</i>	The <a href="#">Particle</a> object to insert or modify in the event
--------------	--

#### Returns

- 1 if a new [Particle](#) object has been inserted in the event
- 0 if an existing [Particle](#) object has been modified
- -1 if the requested role to edit is undefined or incorrect

##### 4.1.2.4 void Event::Store ( std::ofstream \*, double *weight\_* = 1. )

Stores in a file (raw format) all the kinematics on the outgoing leptons

## Parameters

<code>weight__</code>	The weight of the event
-----------------------	-------------------------

## 4.1.2.5 void Event::StoreLHERRecord ( std::ofstream \* of\_, const double weight\_ = 1. )

Stores in a LHE format (a XML-style) all the information on the particles composing this event

## Parameters

<code>of__</code>	The file stream on which the event record has to be saved
<code>weight__</code>	The weight of the event

The documentation for this class was generated from the following file:

- include/event.h

## 4.2 GamGam Class Reference

Computes the matrix element for a  $\gamma\gamma \rightarrow \ell^+\ell^-$  process.

## Public Member Functions

- [GamGam](#) (const unsigned int ndim\_, int nOpt\_, double x\_[])  
*Class constructor.*
- void [ComputeSqS](#) ()  
*Computes  $\sqrt{s}$  for the system.*
- double [ComputeXsec](#) (int nm\_=1)  
*Computes the process' cross section.*
- [Particle](#) \* [GetParticle](#) (int role\_)  
*Get a particle given its role in the process.*
- double [GetT1](#) ()
- void [GetT1extrema](#) (double &t1min\_, double &t1max\_)
- double [GetT2](#) ()
- void [GetT2extrema](#) (double &t2min\_, double &t2max\_)
- bool [IsKinematicsDefined](#) ()  
*Is the system's kinematics well defined?*
- void [SetCuts](#) ([GamGamKinematics](#) cuts\_)  
*Sets the list of kinematic cuts to apply on the outgoing particles' final state.*
- bool [SetIncomingKinematics](#) (int part\_, double momentum\_[3], int pdgId\_)  
*Sets the momentum and PDG id for the incoming particles.*
- bool [SetIncomingKinematics](#) ([Particle](#) ip1\_, [Particle](#) ip2\_)  
*Sets the momentum and PDG id for the incoming particles.*
- bool [SetOutgoingParticles](#) (int part\_, int pdgId\_)  
*Sets the PDG id for the outgoing particles.*

## 4.2.1 Detailed Description

Full class of methods and objects to compute the full analytic matrix element [2] for the  $\gamma\gamma \rightarrow \ell^+\ell^-$  process according to a set of kinematic constraints provided for the incoming and outgoing particles (the [GamGamKinematics](#) object).

## 4.2.2 Constructor &amp; Destructor Documentation

4.2.2.1 GamGam::GamGam ( const unsigned int *ndim\_*, int *nOpt\_*, double *x\_[]* )

Sets the mandatory parameters used in the methods computing the kinematics and the cross-section of this phase space point.

## Parameters

<i>ndim_</i>	The number of dimensions of the point in the phase space
<i>nOpt_</i>	Optimisation???
<i>x_[]</i>	The ( <i>ndim_</i> )-dimensional point in the phase space on which the kinematics and the cross-section are computed

**Todo** Figure out how this *nOpt\_* parameter is affecting the final cross-section computation and events generation

## 4.2.3 Member Function Documentation

## 4.2.3.1 void GamGam::ComputeSqS ( )

Computes the centre of mass energy for the system, according to the incoming particles' kinematics

4.2.3.2 double GamGam::ComputeXsec ( int *nm\_* = 1 )

Computes the cross-section for the  $\gamma\gamma \rightarrow \ell^+\ell^-$  process with the given kinematics

## Returns

$\frac{d\sigma}{dx}(\gamma\gamma \rightarrow \ell^+\ell^-)$ , the differential cross-section for the given point in the phase space.

4.2.3.3 Particle\* GamGam::GetParticle ( int *role\_* )

## Parameters

<i>role_</i>	An integer denoting the particle's role in the selected production process
--------------	--

## 4.2.3.4 double GamGam::GetT1 ( ) [inline]

Returns the value for the first photon virtuality

## Returns

$t_1$ , the first photon virtuality

4.2.3.5 void GamGam::GetT1extrema ( double & *t1min\_*, double & *t1max\_* ) [inline]

Returns the two limit values for the first photon virtuality

## Parameters

<i>t1min_</i>	The minimal value for $t_1$
<i>t1max_</i>	The maximal value for $t_1$

## 4.2.3.6 double GamGam::GetT2 ( ) [inline]

Returns the value for the second photon virtuality

## Returns

$t_2$ , the second photon virtuality

**4.2.3.7** `void GamGam::GetT2extrema ( double & t2min_, double & t2max_ ) [inline]`

Returns the two limit values for the second photon virtuality

## Parameters

<code>t2min_</code>	The minimal value for $t_2$
<code>t2max_</code>	The maximal value for $t_2$

**4.2.3.8** `bool GamGam::IsKinematicsDefined ( ) [inline]`

Is the system's kinematics well defined and compatible with the process ? This check is mandatory to perform the (`_ndim`)-dimensional point's cross-section computation.

## Returns

A boolean stating if the input kinematics and the final states are well defined

**4.2.3.9** `void GamGam::SetCuts ( GamGamKinematics cuts_ )`

## Parameters

<code>cuts_</code>	The Cuts object containing the kinematic parameters
--------------------	---

**4.2.3.10** `bool GamGam::SetIncomingKinematics ( int part_, double momentum_[3], int pdgld_ )`

Specifies the incoming particles' kinematics as well as their properties (role in the process and code according to the PDG convention)

## Parameters

<code>part_</code>	Role of the particle in the process
<code>momentum_[ ]</code>	3-momentum of the particle
<code>pdgld_</code>	<a href="#">Particle</a> ID according to the PDG convention

## Returns

True if the kinematics was correctly set for the given particle role

**4.2.3.11** `bool GamGam::SetIncomingKinematics ( Particle ip1_, Particle ip2_ )`

Specifies the incoming particles' kinematics as well as their properties using two [Particle](#) objects.

## Parameters

<code>ip1_</code>	Information on the first incoming particle
<code>ip2_</code>	Information on the second incoming particle

**4.2.3.12** `bool GamGam::SetOutgoingParticles ( int part_, int pdgld_ )`

## Parameters

<code>part_</code>	Role of the particle in the process
<code>pdgld_</code>	<a href="#">Particle</a> ID according to the PDG convention

The documentation for this class was generated from the following file:

- include/gamgam.h

### 4.3 GamGamKinematics Class Reference

List of kinematic cuts to apply on the central and outgoing phase space.

#### Data Fields

- double `emax`  
*Maximal energy of the central two-photons system.*
- double `emin`  
*Minimal energy of the central two-photons system.*
- int `kinematics`  
*Type of kinematics to consider for the phase space.*
- int `mode`  
*Sets of cuts to apply on the final phase space.*
- double `ptmax`  
*Maximal transverse momentum of the single outgoing leptons.*
- double `ptmin`  
*Minimal transverse momentum of the single outgoing leptons.*
- double `q2max`  
*The maximal value of  $Q^2$ .*
- double `q2min`  
*The minimal value of  $Q^2$ .*
- double `thetamax`  
*Maximal polar ( $\theta_{\max}$ ) angle of the outgoing leptons, expressed in degrees.*
- double `thetamin`  
*Minimal polar ( $\theta_{\min}$ ) angle of the outgoing leptons, expressed in degrees.*
- double `wmax`  
*The maximal  $s$  on which the cross section is integrated. If negative, the maximal energy available to the system (hence,  $s = (\sqrt{s})^2$ ) is provided.*
- double `wmin`  
*The minimal  $s$  on which the cross section is integrated.*

#### 4.3.1 Field Documentation

##### 4.3.1.1 int GamGamKinematics::kinematics

Type of kinematics to consider for the process. Can either be :

- 0 for the electron-electron elastic case
- 1 for the proton-proton elastic case
- 2 for the proton-proton single-dissociative (or inelastic) case
- 3 for the proton-proton double-dissociative case

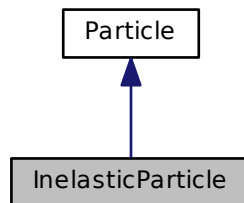
The documentation for this class was generated from the following file:

- include/gamgam.h



## 4.4 InelasticParticle Class Reference

Inheritance diagram for InelasticParticle:



### Public Member Functions

- void [AddDaughter](#) ([Particle](#) \*part\_)  
*Specify a decay product for this particle.*
- void [E](#) (double E\_)  
*Sets the particle's energy.*
- [Particle](#) \* [GetDaughter](#) (const unsigned int num\_=0)  
*Gets a daughter from this particle, labelled by its identifier in this particle's daughters list.*
- std::string [GetLHEline](#) (bool revert\_=false)
- [Particle](#) \* [GetMother](#) ()  
*Gets the mother particle from which this particle arises.*
- void [Hadronise](#) ()  
*Hadronises the particle using Pythia.*
- double [M](#) ()  
*Gets the particle's mass.*
- double [M2](#) ()  
*Gets the particle's squared mass.*
- unsigned int [NumDaughters](#) ()  
*Gets the number of daughter particles arising from this one.*
- bool [P](#) (double px\_, double py\_, double pz\_)  
*Sets the 3-momentum associated to the particle.*
- bool [P](#) (double px\_, double py\_, double pz\_, double E\_)  
*Sets the 4-momentum associated to the particle.*
- bool [P](#) (double p\_[3], double E\_)  
*Sets the 4-momentum associated to the particle.*
- void [SetMother](#) ([Particle](#) \*part\_)  
*Sets the mother particle (from which this particle arises)*

### Data Fields

- double [eta](#)  
*Pseudo-rapidity.*
- bool [isValid](#)  
*Is this particle a valid particle which can be used for kinematic computations ?*

- double [p](#)  
*Norm of the 3-momentum, in GeV/c.*
- int [pdgId](#)  
*Particle Data Group integer identifier.*
- double [pt](#)  
*Transverse momentum, in GeV/c.*
- double [px](#)  
*Momentum along the x-axis in GeV/c.*
- double [py](#)  
*Momentum along the y-axis in GeV/c.*
- double [pz](#)  
*Momentum along the z-axis in GeV/c.*
- int [role](#)  
*Role in the considered process.*
- int [status](#)  
*Particle status.*

#### 4.4.1 Detailed Description

Class containing the information on a particle supposed to decay or fragment in the process

#### 4.4.2 Member Function Documentation

##### 4.4.2.1 void Particle::AddDaughter ( Particle \* *part\_* ) [inherited]

Parameters

<i>part_</i>	The <a href="#">Particle</a> object in which this particle will desintegrate or convert
--------------	---

##### 4.4.2.2 void Particle::E ( double *E\_* ) [inline], [inherited]

Parameters

<i>E_</i>	Energy, in GeV
-----------	----------------

##### 4.4.2.3 Particle\* Particle::GetDaughter ( const unsigned int *num\_* = 0 ) [inherited]

Returns

A [Particle](#) object containing all the kinematic information related to this daughter particle

##### 4.4.2.4 std::string Particle::GetLHEline ( bool *revert\_* = false ) [inherited]

Returns a string containing all the particle's kinematics as expressed in the Les Houches format

Parameters

<i>revert_</i>	Is the event symmetric ? If set to true, the third component of the momentum is reverted.
----------------	---

**Returns**

The LHE line associated to the particle, and containing the particle's history (mother/daughters), its kinematics, and its status

**4.4.2.5 void InelasticParticle::Hadronise ( )**

Hadronises the particle with Pythia, and builds the shower (list of [Particle](#) objects) embedded in this object

**4.4.2.6 double Particle::M ( ) [inline], [inherited]**

Gets the particle's mass in  $\text{GeV}/c^2$ .

**Returns**

The particle's mass

**4.4.2.7 bool Particle::P ( double *px\_*, double *py\_*, double *pz\_* ) [inline], [inherited]****Parameters**

<i>px_</i>	Momentum along the <i>x</i> -axis, in $\text{GeV}/c$
<i>py_</i>	Momentum along the <i>y</i> -axis, in $\text{GeV}/c$
<i>pz_</i>	Momentum along the <i>z</i> -axis, in $\text{GeV}/c$

Here is the call graph for this function:

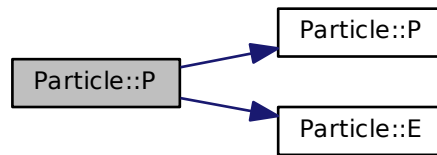
**4.4.2.8 bool Particle::P ( double *px\_*, double *py\_*, double *pz\_*, double *E\_* ) [inline], [inherited]**

Sets the 4-momentum associated to the particle, and computes its (invariant) mass.

**Parameters**

<i>px_</i>	Momentum along the <i>x</i> -axis, in $\text{GeV}/c$
<i>py_</i>	Momentum along the <i>y</i> -axis, in $\text{GeV}/c$
<i>pz_</i>	Momentum along the <i>z</i> -axis, in $\text{GeV}/c$
<i>E_</i>	Energy, in $\text{GeV}$

Here is the call graph for this function:



4.4.2.9 `bool Particle::P ( double  $p_{\perp}$ , double  $E_{\perp}$  )` [inherited]

Parameters

$p_{\perp}$	3-momentum
$E_{\perp}$	Energy, in GeV

4.4.2.10 `void Particle::SetMother ( Particle *  $part_{\perp}$  )` [inline], [inherited]

Parameters

$part_{\perp}$	A <a href="#">Particle</a> object containing all the information on the mother particle
----------------	---

### 4.4.3 Field Documentation

4.4.3.1 `int Particle::status` [inherited]

Codes 1-10 correspond to currently existing partons/particles, and larger codes contain partons/particles which no longer exist, or other kinds of event information

The documentation for this class was generated from the following file:

- `include/inelastic.h`

## 4.5 InputParameters Class Reference

List of input parameters used to start and run the simulation job.

### Public Member Functions

- `void Dump ()`  
*Dumps the input parameters in the console.*
- `bool ReadConfigFile (std::string inFile_)`  
*Reads content from config file to load the variables.*
- `void SetEtaRange (double etamin_, double etamax_)`  
*Sets the pseudo-rapidity range for the produced leptons.*
- `bool StoreConfigFile (std::string outFile_)`  
*Stores the full run configuration to an external config file.*

## Data Fields

- bool [debug](#)  
*Do we need control plots all along the process?*
- std::ofstream \* [file](#)  
*The file in which to store the events generation's output.*
- bool [generation](#)  
*Are we generating events ? (true) or are we only computing the cross-section ? (false)*
- int [gpdf](#)  
*PDFLIB group to use.*
- double [in1p](#)  
*First incoming particle's momentum (in GeV/c)*
- double [in2p](#)  
*Second incoming particle's momentum (in GeV/c)*
- int [itvg](#)  
*Maximal number of iterations to perform by VEGAS.*
- double [maxenergy](#)  
*Maximal energy of the outgoing leptons.*
- int [maxgen](#)  
*Maximal number of events to generate in this run.*
- double [maxmx](#)  
*Maximal  $M_X$  of the outgoing proton remnants.*
- double [maxpt](#)  
*Maximal  $p_T$  of the outgoing leptons.*
- double [maxtheta](#)  
*Maximal polar angle  $\theta$  of the outgoing leptons.*
- int [mcut](#)  
*Set of cuts to apply on the outgoing leptons.*
- double [minenergy](#)  
*Minimal energy of the outgoing leptons.*
- double [minmx](#)  
*Minimal  $M_X$  of the outgoing proton remnants.*
- double [minpt](#)  
*Minimal  $p_T$  of the outgoing leptons.*
- double [mintheta](#)  
*Minimal polar angle  $\theta$  of the outgoing leptons.*
- int [ngen](#)  
*Number of events already generated in this run.*
- int [ntreat](#)  
*Maximal number of TREAT calls.*
- int [p1mod](#)  
*First particle's mode.*
- int [p2mod](#)  
*Second particle's mode.*
- int [pair](#)  
*PDG id of the outgoing leptons.*
- int [qpdf](#)  
*Number of quarks.*
- int [spdf](#)  
*PDFLIB set to use.*
- bool [store](#)  
*Are the events generated in this run to be stored in the output file ?*
- bool [symmetrise](#)  
*Control plots objects.*

## 4.5.1 Detailed Description

## Note

The default parameters are derived from GMUINI in LPAIR

## 4.5.2 Member Function Documentation

## 4.5.2.1 bool InputParameters::ReadConfigFile ( std::string inFile\_ )

Reads the list of parameters to be used in this cross-section computation/events generation from an external input card.

## Parameters

<i>inFile_</i>	Name of the configuration file to load
----------------	--

## 4.5.2.2 void InputParameters::SetEtaRange ( double etamin\_, double etamax\_ )

Defines the range to cover in pseudo-rapidity for the outgoing leptons produced in this process. This method converts this range into a range in  $\theta$ , the polar angle.

## Parameters

<i>etamin_</i>	The minimal value of $\eta$ for the outgoing leptons
<i>etamax_</i>	The maximal value of $\eta$ for the outgoing leptons

## 4.5.2.3 bool InputParameters::StoreConfigFile ( std::string outFile\_ )

## Parameters

<i>outFile_</i>	Name of the configuration file to create
-----------------	--

## 4.5.3 Field Documentation

## 4.5.3.1 bool InputParameters::debug

Enables or disables the production of control plots for several kinematic quantities in this process

## 4.5.3.2 double InputParameters::maxmx

Maximal mass of the outgoing proton remnants,  $M_X$ , in  $\text{GeV}/c^2$ .

## 4.5.3.3 double InputParameters::maxpt

Maximal transverse momentum cut to apply on the outgoing lepton(s)

## 4.5.3.4 int InputParameters::mcut

Set of cuts to apply on the outgoing leptons in order to restrain the available kinematic phase space :

- 0 - No cuts at all (for the total cross section)
- 1 - Vermaserens' hypothetical detector cuts : for both leptons,
  - $\frac{|p_z|}{|\mathbf{p}|} \leq 0.75$  and  $p_T \geq 1 \text{ GeV}/c$ , or
  - $0.75 < \frac{|p_z|}{|\mathbf{p}|} \leq 0.95$  and  $p_z > 1 \text{ GeV}/c$ ,
- 2 - Cuts on both the outgoing leptons, according to the provided cuts parameters

- 3 - Cuts on at least one outgoing lepton, according to the provided cut parameters

#### 4.5.3.5 double InputParameters::minmx

Minimal mass of the outgoing proton remnants,  $M_X$ , in  $\text{GeV}/c^2$ .

#### 4.5.3.6 double InputParameters::minpt

Minimal transverse momentum cut to apply on the outgoing lepton(s)

#### 4.5.3.7 int InputParameters::ntreat

##### Note

Is it correctly implemented ?

#### 4.5.3.8 int InputParameters::p1mod

The first incoming particle type and kind of interaction :

- 1 - electron,
- 2 - proton elastic,
- 3 - proton inelastic without parton treatment,
- 4 - proton inelastic in parton model

##### Note

Was named PMOD in ILPAIR

#### 4.5.3.9 int InputParameters::p2mod

##### Note

Was named EMOD in ILPAIR

#### 4.5.3.10 int InputParameters::pair

The particle code of produced leptons, as defined by the PDG convention :

- 11 - for  $e^+e^-$  pairs
- 13 - for  $\mu^+\mu^-$  pairs
- 15 - for  $\tau^+\tau^-$  pairs

#### 4.5.3.11 bool InputParameters::symmetrise

List of Gnuplot objects which can be used to produce control plots all along the cross-section determination and events generation process

##### Note

Maximum number of these can be raised in the [utils.h](#) file, but pay attention to the memory load since these Gnuplot objects are still under development!

The documentation for this class was generated from the following file:

- include/utils.h

## 4.6 MCGen Class Reference

Core of the Monte-Carlo generator.

### Public Member Functions

- [MCGen](#) ([InputParameters](#) ip\_)
- Class constructor.*
- void [ComputeXsection](#) (double \*, double \*)
- [InputParameters](#) [GetInputParameters](#) ()
- Returns the set of parameters used to setup the phase space to integrate.*

### 4.6.1 Detailed Description

This object represents the core of this Monte Carlo generator, with its allowance to generate the events (using the embedded [Vegas](#) object) and to study the phase space in term of the variation of resulting cross section while scanning the various parameters (point  $\mathbf{x}$  in the DIM-dimensional phase space).

The phase space is constrained using the [InputParameters](#) object given as an argument to the constructor, and the differential cross-sections for each value of the array  $\mathbf{x}$  are computed in the f-function defined outside (but populated inside) this object.

This f-function embeds a [GamGam](#) object which defines all the methods to obtain this differential cross-section as well as the in- and outgoing kinematics associated to each particle.

### Author

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### Date

February 2013

### 4.6.2 Constructor & Destructor Documentation

#### 4.6.2.1 MCGen::MCGen ( InputParameters ip\_ )

Sets the number of dimensions on which to perform the integration, according to the set of input parameters given as an argument and propagated to the whole object

#### Parameters

<i>ip_</i>	List of input parameters defining the phase space on which to perform the integration
------------	---

### 4.6.3 Member Function Documentation

#### 4.6.3.1 void MCGen::ComputeXsection ( double \*, double \* )

Computes the cross-section for the run defined by this object. This returns the cross-section as well as the absolute error computed along.

#### 4.6.3.2 InputParameters MCGen::GetInputParameters ( ) [inline]

#### Returns

The InputParameter object embedded in this class

The documentation for this class was generated from the following file:

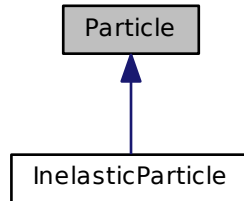


- `include/mcgen.h`

## 4.7 Particle Class Reference

Kinematics of one particle.

Inheritance diagram for Particle:



### Public Member Functions

- `void AddDaughter (Particle *part_)`  
*Specify a decay product for this particle.*
- `void E (double E_)`  
*Sets the particle's energy.*
- `Particle * GetDaughter (const unsigned int num_=0)`  
*Gets a daughter from this particle, labelled by its identifier in this particle's daughters list.*
- `std::string GetLHEline (bool revert_=false)`
- `Particle * GetMother ()`  
*Gets the mother particle from which this particle arises.*
- `double M ()`  
*Gets the particle's mass.*
- `double M2 ()`  
*Gets the particle's squared mass.*
- `unsigned int NumDaughters ()`  
*Gets the number of daughter particles arising from this one.*
- `bool P (double px_, double py_, double pz_)`  
*Sets the 3-momentum associated to the particle.*
- `bool P (double px_, double py_, double pz_, double E_)`  
*Sets the 4-momentum associated to the particle.*
- `bool P (double p_[3], double E_)`  
*Sets the 4-momentum associated to the particle.*
- `void SetMother (Particle *part_)`  
*Sets the mother particle (from which this particle arises)*

## Data Fields

- double [eta](#)  
*Pseudo-rapidity.*
- bool [isValid](#)  
*Is this particle a valid particle which can be used for kinematic computations ?*
- double [p](#)  
*Norm of the 3-momentum, in GeV/c.*
- int [pdgId](#)  
*Particle Data Group integer identifier.*
- double [pt](#)  
*Transverse momentum, in GeV/c.*
- double [px](#)  
*Momentum along the x-axis in GeV/c.*
- double [py](#)  
*Momentum along the y-axis in GeV/c.*
- double [pz](#)  
*Momentum along the z-axis in GeV/c.*
- int [role](#)  
*Role in the considered process.*
- int [status](#)  
*Particle status.*

## 4.7.1 Detailed Description

Kinematic information for one particle

## 4.7.2 Member Function Documentation

4.7.2.1 void Particle::AddDaughter ( Particle \* *part\_* )

## Parameters

<i>part_</i>	The <a href="#">Particle</a> object in which this particle will desintegrate or convert
--------------	---

4.7.2.2 void Particle::E ( double *E\_* ) [inline]

## Parameters

<i>E_</i>	Energy, in GeV
-----------	----------------

4.7.2.3 Particle\* Particle::GetDaughter ( const unsigned int *num\_* = 0 )

## Returns

A [Particle](#) object containing all the kinematic information related to this daughter particle

4.7.2.4 std::string Particle::GetLHEline ( bool *revert\_* = false )

Returns a string containing all the particle's kinematics as expressed in the Les Houches format

## Parameters

<i>revert_</i>	Is the event symmetric ? If set to true, the third component of the momentum is reverted.
----------------	---

**Returns**

The LHE line associated to the particle, and containing the particle's history (mother/daughters), its kinematics, and its status

**4.7.2.5 double Particle::M ( ) [inline]**

Gets the particle's mass in  $\text{GeV}/c^2$ .

**Returns**

The particle's mass

**4.7.2.6 bool Particle::P ( double *px\_*, double *py\_*, double *pz\_* ) [inline]****Parameters**

<i>px_</i>	Momentum along the <i>x</i> -axis, in $\text{GeV}/c$
<i>py_</i>	Momentum along the <i>y</i> -axis, in $\text{GeV}/c$
<i>pz_</i>	Momentum along the <i>z</i> -axis, in $\text{GeV}/c$

Here is the call graph for this function:

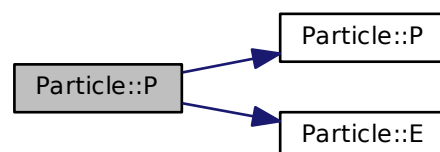
**4.7.2.7 bool Particle::P ( double *px\_*, double *py\_*, double *pz\_*, double *E\_* ) [inline]**

Sets the 4-momentum associated to the particle, and computes its (invariant) mass.

**Parameters**

<i>px_</i>	Momentum along the <i>x</i> -axis, in $\text{GeV}/c$
<i>py_</i>	Momentum along the <i>y</i> -axis, in $\text{GeV}/c$
<i>pz_</i>	Momentum along the <i>z</i> -axis, in $\text{GeV}/c$
<i>E_</i>	Energy, in $\text{GeV}$

Here is the call graph for this function:



4.7.2.8 `bool Particle::P ( double  $p_{-}[3]$ , double  $E_{-}$  )`

## Parameters

$p_{-}$	3-momentum
$E_{-}$	Energy, in GeV

4.7.2.9 `void Particle::SetMother ( Particle *  $part_{-}$  ) [inline]`

## Parameters

$part_{-}$	A <a href="#">Particle</a> object containing all the information on the mother particle
------------	---

## 4.7.3 Field Documentation

4.7.3.1 `int Particle::status`

Codes 1-10 correspond to currently existing partons/particles, and larger codes contain partons/particles which no longer exist, or other kinds of event information

The documentation for this class was generated from the following file:

- `include/particle.h`

## 4.8 Vegas Class Reference

[Vegas](#) Monte-Carlo integrator instance.

## Public Member Functions

- [Vegas](#) (const int  $dim_{-}$ , double  $f_{-}$ (double \*, size\_t, void \*), [InputParameters](#) \* $inParam_{-}$ )
- [~Vegas](#) ()  
*Class destructor.*
- int [Integrate](#) (double \* $result_{-}$ , double \* $abserr_{-}$ )  
*Launches the integration of the provided function.*
- int [LaunchGeneration](#) ()  
*Launches the generation of events.*
- int [MyIntegrate](#) (double \* $result_{-}$ , double \* $abserr_{-}$ )

## 4.8.1 Constructor &amp; Destructor Documentation

4.8.1.1 `Vegas::Vegas ( const int  $dim_{-}$ , double  $f_{-}$ double *, size_t, void *, InputParameters *  $inParam_{-}$  )`

Constructs the class by booking the memory and structures for the GSL [Vegas](#) integrator. This code from the GNU scientific library is based on the [Vegas](#) Monte Carlo integration algorithm developed by P. Lepage. [1]

## Parameters

$dim_{-}$	The number of dimensions on which the function will be integrated
$f_{-}$	The function one is required to integrate
$inParam_{-}$	A list of parameters to define the phase space on which this integration is performed (embedded in an <a href="#">InputParameters</a> object)

## 4.8.2 Member Function Documentation

#### 4.8.2.1 int Vegas::Integrate ( double \* *result\_*, double \* *abserr\_* )

Launches the [Vegas](#) integration of the provided function with the provided input parameters.

##### Parameters

<i>result_</i>	The cross section as integrated by <a href="#">Vegas</a> for the given phase space restrictions
<i>abserr_</i>	The error associated to the computed cross section

#### 4.8.2.2 int Vegas::LaunchGeneration ( )

Launches the [Vegas](#) generation of events according to the provided input parameters.

#### 4.8.2.3 int Vegas::MyIntegrate ( double \* *result\_*, double \* *abserr\_* )

[Vegas](#) algorithm to perform the (*\_dim*)-dimensional Monte Carlo integration of a given function as described in[1]

##### Author

Primary author : G.P. Lepage  
This C++ implementation : L. Forthomme

##### Date

September 1976  
Reviewed in Apr 1978  
FTN5 version 21 Aug 1984  
This C++ implementation is from 12 Dec 2013

##### Parameters

<i>result_</i>	The cross section as integrated by <a href="#">Vegas</a> for the given phase space restrictions
<i>abserr_</i>	The error associated to the computed cross section

The documentation for this class was generated from the following file:

- include/vegas.h

## References

- [1] G Peter Lepage. A new algorithm for adaptive multidimensional integration. *Journal of Computational Physics*, 27(2):192 – 203, 1978.
- [2] J.A.M. Vermaseren. Two-photon processes at very high energies. *Nuclear Physics B*, 229(2):347 – 371, 1983.

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