

LPAIR++  
0.1

Generated by Doxygen 1.8.3.1

Tue Jul 16 2013 15:15:00

## Contents

<b>1</b>	<b>Todo List</b>	<b>1</b>
<b>2</b>	<b>Data Structure Index</b>	<b>1</b>
2.1	Data Structures . . . . .	2
<b>3</b>	<b>Data Structure Documentation</b>	<b>2</b>
3.1	Cuts Class Reference . . . . .	2
3.1.1	Constructor & Destructor Documentation . . . . .	3
3.1.2	Field Documentation . . . . .	3
3.2	GamGam Class Reference . . . . .	3
3.2.1	Constructor & Destructor Documentation . . . . .	4
3.2.2	Member Function Documentation . . . . .	4
3.3	InputParameters Class Reference . . . . .	6
3.3.1	Detailed Description . . . . .	7
3.3.2	Constructor & Destructor Documentation . . . . .	7
3.3.3	Member Function Documentation . . . . .	7
3.3.4	Field Documentation . . . . .	7
3.4	MCGen Class Reference . . . . .	9
3.4.1	Detailed Description . . . . .	9
3.4.2	Constructor & Destructor Documentation . . . . .	9
3.4.3	Member Function Documentation . . . . .	10
3.5	Particle Class Reference . . . . .	11
3.5.1	Detailed Description . . . . .	11
3.5.2	Constructor & Destructor Documentation . . . . .	11
3.5.3	Member Function Documentation . . . . .	11
3.5.4	Field Documentation . . . . .	12
3.6	Vegas Class Reference . . . . .	12
3.6.1	Constructor & Destructor Documentation . . . . .	12
3.6.2	Member Function Documentation . . . . .	13
	<b>Index</b>	<b>14</b>

## 1 Todo List

### Global **GamGam::GamGam** (const unsigned int, double, double, int, double x\_[])

Figure out how this nOpt\_ parameter is affecting the final cross-section computation and events generation  
 What are these w12, w31, w52 parameters introduced in the GAMGAM subroutine ? And why are they set to 0. ?

## 2 Data Structure Index

### 2.1 Data Structures

Here are the data structures with brief descriptions:

<b>Cuts</b>	
List of kinematic cuts to apply on the central and outgoing phase space	2
<b>GamGam</b>	
Computes the matrix element for a $\gamma\gamma \rightarrow \ell^+\ell^-$ process	3
<b>InputParameters</b>	
List of input parameters used to start and run the simulation job	6
<b>MCGen</b>	
Core of the Monte-Carlo generator ; Computes the cross section for any value of the input parameters by calling <b>Vegas</b> on <b>GamGam</b> objects	9
<b>Particle</b>	
Kinematics of one particle	11
<b>Vegas</b>	
<b>Vegas</b> Monte-Carlo integrator instance	12

## 3 Data Structure Documentation

### 3.1 Cuts Class Reference

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

#### Public Member Functions

- **Cuts** ()
- **~Cuts** ()

#### Data Fields

- double **emax**  
*Maximal energy of the central two-photons system.*
- double **emin**  
*Minimal energy of the central two-photons system.*
- int **mode**  
*Sets of cuts to apply on the final phase space.*
- double **mxmax**
- double **mxmin**
- double **ptmax**  
*Maximal transverse momentum of the single outgoing leptons.*
- double **ptmin**  
*Minimal transverse momentum of the single outgoing leptons.*
- 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.*

## 3.1.1 Constructor &amp; Destructor Documentation

3.1.1.1 Cuts::Cuts ( )

3.1.1.2 Cuts::~~Cuts ( )

## 3.1.2 Field Documentation

3.1.2.1 double Cuts::emax

3.1.2.2 double Cuts::emin

3.1.2.3 int Cuts::mode

3.1.2.4 double Cuts::mxmax

3.1.2.5 double Cuts::mxmin

3.1.2.6 double Cuts::ptmax

3.1.2.7 double Cuts::ptmin

3.1.2.8 double Cuts::thetamax

3.1.2.9 double Cuts::thetamin

## 3.2 GamGam Class Reference

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

## Public Member Functions

- [GamGam](#) (const unsigned int, double, double, int, double x\_<sub>0</sub>[])  
*Class constructor.*
- [~GamGam](#) ( )
- void [ComputeSqS](#) ( )  
*Computes  $\sqrt{s}$  for the system.*
- double [ComputeXsec](#) (int nm\_<sub>0</sub>=1)  
*Computes the process' cross section.*
- void [FillKinematics](#) ( )
- [Particle](#) [GetParticle](#) (int)  
*Get a particle given its role in the process.*
- bool [IsKinematicsDefined](#) ( )  
*Is the system's kinematics well defined?*
- void [SetCuts](#) (Cuts)  
*Sets the list of kinematic cuts to apply on the outgoing particles' final state.*
- bool [SetIncomingKinematics](#) (int, double[], int)  
*Sets the momentum and PDG id for the incoming particles.*
- bool [SetOutgoingParticles](#) (int, int)  
*Sets the PDG id for the outgoing particles.*
- void [SetWRange](#) (double, double)  
*Sets the energy range available for the phase space integration.*
- void [StoreEvent](#) (std::ofstream \*, double)

## 3.2.1 Constructor &amp; Destructor Documentation

3.2.1.1 GamGam::GamGam ( const unsigned int *ndim\_*, double *q2min\_*, double *q2max\_*, 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>q2min_</i>	The minimal value of $Q^2$
<i>q2max_</i>	The maximal value of $Q^2$
<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

What are these *w12*, *w31*, *w52* parameters introduced in the GAMGAM subroutine ? And why are they set to 0. ?

## 3.2.1.2 GamGam::~GamGam ( )

## 3.2.2 Member Function Documentation

## 3.2.2.1 void GamGam::ComputeSqS ( )

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

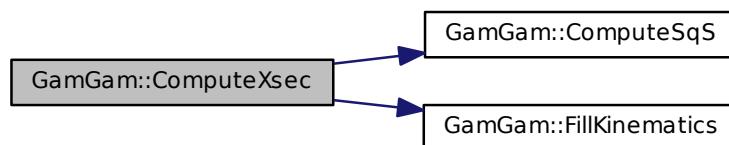
3.2.2.2 double GamGam::ComputeXsec ( int *nm\_* = 1 )

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.

Here is the call graph for this function:



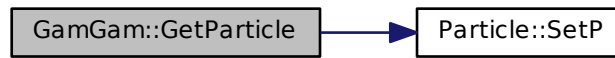
## 3.2.2.3 void GamGam::FillKinematics ( )

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

## Parameters

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

Here is the call graph for this function:



### 3.2.2.5 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

### 3.2.2.6 void GamGam::SetCuts ( Cuts cuts\_ )

#### Parameters

<i>cuts_</i>	The <a href="#">Cuts</a> object containing the kinematic parameters
--------------	---

### 3.2.2.7 bool GamGam::SetIncomingKinematics ( int, double [], int )

Specifies the incoming particles' kinematics as well as their properties (role in the process and PDG Id)

#### Parameters

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

#### Returns

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

### 3.2.2.8 bool GamGam::SetOutgoingParticles ( int *part\_*, int *pdgId\_* )

#### Parameters

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

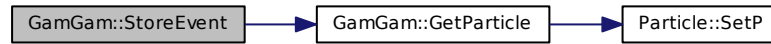
### 3.2.2.9 void GamGam::SetWRange ( double *wmin\_*, double *wmax\_* )

#### Parameters

<i>wmin_</i>	The minimal $s$ on which the cross section is integrated
<i>wmax_</i>	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.

## 3.2.2.10 void GamGam::StoreEvent ( std::ofstream \* file, double weight )

Here is the call graph for this function:



## 3.3 InputParameters Class Reference

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

## Public Member Functions

- [InputParameters](#) ()
- [~InputParameters](#) ()
- bool [ReadConfigFile](#) (std::string)  
*Reads content from config file to load the variables.*
- bool [StoreConfigFile](#) (std::string)  
*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)*
- double [in1p](#)  
*First incoming particle's momentum (in GeV/c)*
- double [in2p](#)  
*Second incoming particle's momentum (in GeV/c)*
- int [itvg](#)  
*Number of [Vegas](#) integrations.*
- double [maxenergy](#)
- double [maxmx](#)
- double [maxpt](#)  
*Maximal transverse momentum of the outgoing leptons.*
- double [maxtheta](#)
- int [mcut](#)  
*Set of cuts to apply on the outgoing leptons.*
- double [minenergy](#)
- double [minmx](#)
- double [minpt](#)  
*Minimal transverse momentum of the outgoing leptons.*
- double [mintheta](#)
- int [ncvg](#)

- int `ngen`
- int `p1mod`  
*First particle's mode.*
- int `p2mod`  
*Second particle's mode.*
- int `pair`  
*PDG id of the outgoing leptons.*
- Gnuplot \* `plot` [MAX\_HISTOS]  
*Control plots objects.*
- bool `store`

### 3.3.1 Detailed Description

#### Note

The default parameters are derived from GMUINI in LPAIR

### 3.3.2 Constructor & Destructor Documentation

#### 3.3.2.1 InputParameters::InputParameters ( )

#### 3.3.2.2 InputParameters::~~InputParameters ( )

### 3.3.3 Member Function Documentation

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

#### Parameters

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

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

#### Parameters

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

### 3.3.4 Field Documentation

#### 3.3.4.1 bool InputParameters::debug

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

#### 3.3.4.2 std::ofstream\* InputParameters::file

#### 3.3.4.3 bool InputParameters::generation

#### 3.3.4.4 double InputParameters::in1p

#### 3.3.4.5 double InputParameters::in2p

#### 3.3.4.6 int InputParameters::itvg

#### 3.3.4.7 double InputParameters::maxenergy

#### 3.3.4.8 double InputParameters::maxmx

#### 3.3.4.9 double InputParameters::maxpt



3.3.4.10 double InputParameters::maxtheta

3.3.4.11 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  $p_T \geq 1$  GeV, or
  - $0.75 < \frac{|p_z|}{|\mathbf{p}|} \leq 0.95$  and  $p_z > 1$  GeV,
- 2 - [Cuts](#) according to the provided parameters

3.3.4.12 double InputParameters::minenergy

3.3.4.13 double InputParameters::minmx

3.3.4.14 double InputParameters::minpt

3.3.4.15 double InputParameters::mintheta

3.3.4.16 int InputParameters::ncvg

3.3.4.17 int InputParameters::ngen

3.3.4.18 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

3.3.4.19 int InputParameters::p2mod

Note

Was named EMOD in ILPAIR

3.3.4.20 int InputParameters::pair

The particle code of produced leptons :

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

## 3.3.4.21 Gnuplot\* InputParameters::plot[MAX\_HISTOS]

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!

## 3.3.4.22 bool InputParameters::store

## 3.4 MCGen Class Reference

Core of the Monte-Carlo generator ; Computes the cross section for any value of the input parameters by calling [Vegas](#) on [GamGam](#) objects.

## Public Member Functions

- [MCGen](#) ([InputParameters](#))  
*Class constructor.*
- [~MCGen](#) ()
- void [AnalyzePhaseSpace](#) (const std::string)
- void [ComputeXsection](#) (double \*, double \*)
- [InputParameters](#) [GetInputParameters](#) ()  
*Returns the set of parameters used to setup the phase space to integrate.*
- void [LaunchGen](#) (const unsigned int)
- void [Test](#) ()

## 3.4.1 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 **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 **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.

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## Date

February 2013

## 3.4.2 Constructor &amp; Destructor Documentation

3.4.2.1 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

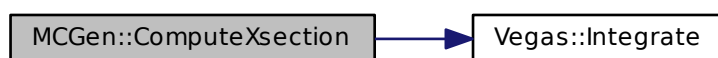
<i>ip_</i>	List of input parameters defining the phase space on which to perform the integration
------------	---

## 3.4.2.2 MCGen::~MCGen ( )

## 3.4.3 Member Function Documentation

3.4.3.1 void MCGen::AnalyzePhaseSpace ( const std::string *outputFile\_* )3.4.3.2 void MCGen::ComputeXsection ( double \* *xsec\_*, double \* *err\_* )

Here is the call graph for this function:



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

## Returns

The InputParameter object embedded in this class

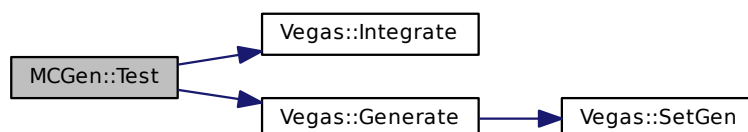
3.4.3.4 void MCGen::LaunchGen ( const unsigned int *count\_* )

Here is the call graph for this function:



## 3.4.3.5 void MCGen::Test ( )

Here is the call graph for this function:



### 3.5 Particle Class Reference

Kinematics of one particle.

#### Public Member Functions

- [Particle](#) ()
- [~Particle](#) ()
- `std::string` [GetLHEline](#) ()
- `void` [SetP](#) (double, double, double)  
*Sets the momentum.*

#### Data Fields

- `double` [e](#)  
*Energy in GeV.*
- `double` [m](#)  
*Mass in  $\text{GeV}/c^2$ .*
- `int` [pdgId](#)  
*Particle Data Group integer identifier.*
- `double` [pt](#)  
*Transverse momentum.*
- `double` [px](#)  
*Momentum along the x-axis in  $\text{GeV}/c$ .*
- `double` [py](#)  
*Momentum along the y-axis in  $\text{GeV}/c$ .*
- `double` [pz](#)  
*Momentum along the z-axis in  $\text{GeV}/c$ .*
- `int` [role](#)  
*Role in the considered process.*

#### 3.5.1 Detailed Description

Kinematic information for one particle

#### 3.5.2 Constructor & Destructor Documentation

##### 3.5.2.1 `Particle::Particle ( )`

##### 3.5.2.2 `Particle::~~Particle ( )`

#### 3.5.3 Member Function Documentation

##### 3.5.3.1 `std::string Particle::GetLHEline ( )`

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

#### Returns

The LHE line

3.5.3.2 void Particle::SetP ( double *px\_*, double *py\_*, double *pz\_* )

## Parameters

<i>px_</i>	Momentum along the <i>x</i> -axis
<i>py_</i>	Momentum along the <i>y</i> -axis
<i>pz_</i>	Momentum along the <i>z</i> -axis

## 3.5.4 Field Documentation

## 3.5.4.1 double Particle::e

## 3.5.4.2 double Particle::m

## 3.5.4.3 int Particle::pdgId

## 3.5.4.4 double Particle::pt

## 3.5.4.5 double Particle::px

## 3.5.4.6 double Particle::py

## 3.5.4.7 double Particle::pz

## 3.5.4.8 int Particle::role

## 3.6 Vegas Class Reference

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

## Public Member Functions

- [Vegas](#) (int, double *f\_*(double \*, size\_t, void \*), [InputParameters](#) \**inParam\_*)
- [~Vegas](#) ()  
*Class destructor.*
- int [Generate](#) (int)
- int [Integrate](#) (double \*, double \*)  
*Launches the integration of the provided function.*
- int [LaunchGeneration](#) (int)  
*Launches the generation of events.*
- void [SetGen](#) ()

## 3.6.1 Constructor &amp; Destructor Documentation

3.6.1.1 Vegas::Vegas ( 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

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

## 3.6.1.2 Vegas::~Vegas ( )

## 3.6.2 Member Function Documentation

3.6.2.1 `int Vegas::Generate ( int nEvt_ )`

Here is the call graph for this function:

3.6.2.2 `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

3.6.2.3 `int Vegas::LaunchGeneration ( int )`

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

## Parameters

<i>nEvts_</i>	The number of events to generate
---------------	----------------------------------

3.6.2.4 `void Vegas::SetGen ( )`

XXXXXX.....

## References

- [1] G Peter Lepage. A new algorithm for adaptive multidimensional integration. *Journal of Computational Physics*, 27(2):192 – 203, 1978.

## Index

- ~Cuts
  - Cuts, [2](#)
- ~GamGam
  - GamGam, [3](#)
- ~InputParameters
  - InputParameters, [6](#)
- ~MCGen
  - MCGen, [9](#)
- ~Particle
  - Particle, [11](#)
- ~Vegas
  - Vegas, [12](#)
- AnalyzePhaseSpace
  - MCGen, [9](#)
- ComputeSqS
  - GamGam, [3](#)
- ComputeXsec
  - GamGam, [3](#)
- ComputeXsection
  - MCGen, [9](#)
- Cuts, [1](#)
  - ~Cuts, [2](#)
  - Cuts, [2](#)
  - emax, [2](#)
  - emin, [2](#)
  - mode, [2](#)
  - mxmax, [2](#)
  - mxmin, [2](#)
  - ptmax, [2](#)
  - ptmin, [2](#)
  - thetamax, [2](#)
  - thetamin, [2](#)
- debug
  - InputParameters, [7](#)
- e
  - Particle, [11](#)
- emax
  - Cuts, [2](#)
- emin
  - Cuts, [2](#)
- file
  - InputParameters, [7](#)
- FillKinematics
  - GamGam, [4](#)
- GamGam, [2](#)
  - ~GamGam, [3](#)
  - ComputeSqS, [3](#)
  - ComputeXsec, [3](#)
  - FillKinematics, [4](#)
  - GamGam, [3](#)
  - GamGam, [3](#)
  - GetParticle, [4](#)
  - IsKinematicsDefined, [4](#)
  - SetCuts, [4](#)
  - SetIncomingKinematics, [4](#)
  - SetOutgoingParticles, [5](#)
  - SetWRange, [5](#)
  - StoreEvent, [5](#)
- Generate
  - Vegas, [12](#)
- generation
  - InputParameters, [7](#)
- GetInputParameters
  - MCGen, [9](#)
- GetLHEline
  - Particle, [11](#)
- GetParticle
  - GamGam, [4](#)
- in1p
  - InputParameters, [7](#)
- in2p
  - InputParameters, [7](#)
- InputParameters, [5](#)
  - ~InputParameters, [6](#)
  - debug, [7](#)
  - file, [7](#)
  - generation, [7](#)
  - in1p, [7](#)
  - in2p, [7](#)
  - InputParameters, [6](#)
  - InputParameters, [6](#)
  - itvg, [7](#)
  - maxenergy, [7](#)
  - maxmx, [7](#)
  - maxpt, [7](#)
  - maxtheta, [7](#)
  - mcut, [7](#)
  - minenergy, [7](#)
  - minmx, [7](#)
  - minpt, [7](#)
  - mintheta, [7](#)
  - ncvg, [7](#)
  - ngen, [7](#)
  - p1mod, [7](#)
  - p2mod, [8](#)
  - pair, [8](#)
  - plot, [8](#)
  - ReadConfigFile, [6](#)
  - store, [8](#)
  - StoreConfigFile, [6](#)
- Integrate
  - Vegas, [12](#)
- IsKinematicsDefined
  - GamGam, [4](#)
- itvg
  - InputParameters, [7](#)



LaunchGen  
  MCGen, 9  
LaunchGeneration  
  Vegas, 12  
  
m  
  Particle, 11  
MCGen, 8  
  ~MCGen, 9  
  AnalyzePhaseSpace, 9  
  ComputeXsection, 9  
  GetInputParameters, 9  
  LaunchGen, 9  
  MCGen, 9  
  MCGen, 9  
  Test, 10  
maxenergy  
  InputParameters, 7  
maxmx  
  InputParameters, 7  
maxpt  
  InputParameters, 7  
maxtheta  
  InputParameters, 7  
mcut  
  InputParameters, 7  
minenergy  
  InputParameters, 7  
minmx  
  InputParameters, 7  
minpt  
  InputParameters, 7  
mintheta  
  InputParameters, 7  
mode  
  Cuts, 2  
mxmax  
  Cuts, 2  
mxmin  
  Cuts, 2  
  
ncvg  
  InputParameters, 7  
ngen  
  InputParameters, 7  
  
p1mod  
  InputParameters, 7  
p2mod  
  InputParameters, 8  
pair  
  InputParameters, 8  
Particle, 10  
  ~Particle, 11  
  e, 11  
  GetLHEline, 11  
  m, 11  
  Particle, 11  
  pdgId, 11  
  pt, 11  
  px, 11  
  py, 11  
  pz, 11  
  role, 11  
  SetP, 11  
pdgId  
  Particle, 11  
plot  
  InputParameters, 8  
pt  
  Particle, 11  
ptmax  
  Cuts, 2  
ptmin  
  Cuts, 2  
px  
  Particle, 11  
py  
  Particle, 11  
pz  
  Particle, 11  
  
ReadConfigFile  
  InputParameters, 6  
role  
  Particle, 11  
  
SetCuts  
  GamGam, 4  
SetGen  
  Vegas, 13  
SetIncomingKinematics  
  GamGam, 4  
SetOutgoingParticles  
  GamGam, 5  
SetP  
  Particle, 11  
SetWRange  
  GamGam, 5  
store  
  InputParameters, 8  
StoreConfigFile  
  InputParameters, 6  
StoreEvent  
  GamGam, 5  
  
Test  
  MCGen, 10  
thetamax  
  Cuts, 2  
thetamin  
  Cuts, 2  
  
Vegas, 11  
  ~Vegas, 12  
  Generate, 12  
  Integrate, 12  
  LaunchGeneration, 12

SetGen, [13](#)  
Vegas, [12](#)