

LPAIR++

0.1

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1 Data Structure Index

1.1 Data Structures

Here are the data structures with brief descriptions:

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2 Data Structure Documentation

2.1 Cuts Class Reference

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

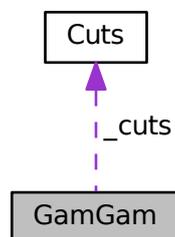
Data Fields

- double `emax`
Maximal energy of the central two-photons system.
- double `emin`
Minimal energy of the central two-photons system.
- int `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 `thetamax`
Maximal polar (θ) angle of the outgoing leptons.
- double `thetamin`
Minimal polar (θ) angle of the outgoing leptons.

2.2 GamGam Class Reference

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

Collaboration diagram for GamGam:



Public Member Functions

- **GamGam** (int, double, double, int, 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)
- bool **IsKinematicsDefined** ()
Is the system's kinematics well defined?
- bool **Orient** ()
Energies/momenta computation for the various particles, in the CM system.
- double **PeriPP** (int, int)
Computes the matrix element squared for the requested process.
- bool **Pickin** ()
- 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.

Private Attributes

- double **_a1**
- double **_a2**
- double **_a5**
- double **_a6**
- double **_acc3**
- double **_acc4**
- double **_al3**
- double **_al4**
- double **_bb**
- double **_be4**
- double **_be5**
- double **_betgam**
 $\beta\gamma$ factor of the centre of mass system, used in the computation of the inverse boost for the outgoing leptons
- double **_cotth1**
- double **_cotth2**
- double **_cp3**
 $\cos\phi_3$ of the first outgoing proton-like particle
- double **_cp5**
 $\cos\phi_5$ of the second outgoing proton-like particle
- double **_cp6**
 $\cos\phi_6$ of the first outgoing lepton
- double **_cp7**
 $\cos\phi_7$ of the second outgoing lepton
- double **_ct3**
 $\cos\theta_3$ of the first outgoing proton-like particle

- double `_ct4`
 $\cos\theta_4$ of the two-photons centre of mass system
- double `_ct5`
 $\cos\theta_5$ of the second outgoing proton-like particle
- double `_ct6`
 $\cos\theta_6$ of the first outgoing lepton
- double `_ct7`
 $\cos\theta_7$ of the second outgoing lepton
- double `_ctcm6`
 $\cos\theta_6^{\text{CM}}$, production angle of the first outgoing lepton, computed in the centre of mass system.
- `Cuts _cuts`
Set of cuts to apply on the final phase space.
- double `_d1`
 $\delta_1 = m_3^2 - m_1^2$ as defined in Vermaseren's paper
- double `_d1dq`
- double `_d1dq2`
- double `_d2`
 $\delta_4 = m_5^2 - m_2^2$ as defined in Vermaseren's paper
- double `_d3`
- double `_d4`
 $\delta_5 = m_4^2 - t_1$ as defined in Vermaseren's paper
- double `_d5`
 $\delta_2 = m_1^2 - m_2^2$ as defined in Vermaseren's paper
- double `_d6`
 $\delta_6 = m_4^2 - m_5^2$ as defined in Vermaseren's paper
- double `_d7`
- double `_d8`
 $\delta_3 = t_1 - m_2^2$ as defined in Vermaseren's paper
- double `_dd1`
- double `_dd2`
- double `_dd3`
- double `_dd4`
- double `_dd5`
- double `_de3`
- double `_de5`
- double `_delta`
- double `_dj`
- double `_e6lab`
 E_6^{lab} , energy of the first outgoing lepton, computed in the lab frame
- double `_e7lab`
 E_7^{lab} , energy of the second outgoing lepton, computed in the lab frame
- double `_ec4`
 E_4 , energy of the two-photon central system
- double `_ecut`
- double `_el6`
 E_6 , energy of the first outgoing lepton
- double `_el7`
 E_7 , energy of the second outgoing lepton
- double `_ep1`
 E_1 , energy of the first proton-like incoming particle
- double `_ep2`
 E_2 , energy of the second proton-like incoming particle

- double `_ep3`
 E_3 , energy of the first proton-like outgoing particle
- double `_ep5`
 E_5 , energy of the second proton-like outgoing particle
- double `_epsi`
- double `_etot`
Total energy provided by the two incoming proton-like particles.
- double `_g4`
- double `_g5`
- double `_g6`
- double `_gamma`
 γ factor of the centre of mass system, used in the computation of the inverse boost for the outgoing leptons
- double `_gram`
- double `_mc4`
 m_4 , mass of the two-photon central system
- double `_ml6`
 m_6 , mass of the first outgoing lepton
- double `_ml7`
 m_7 , mass of the second outgoing lepton
- double `_mp1`
 m_1 , mass of the first proton-like incoming particle
- double `_mp2`
 m_2 , mass of the second proton-like incoming particle
- double `_mp3`
 m_3 , mass of the first proton-like outgoing particle
- double `_mp5`
 m_5 , mass of the second proton-like outgoing particle
- int `_ndim`
Number of dimensions on which the integration has to be performed.
- int `_nOpt`
- double `_p`
- double `_p12`
- double `_p13`
- double `_p14`
- double `_p15`
- double `_p1k2`
- double `_p23`
- double `_p24`
- double `_p25`
- double `_p2k1`
- double `_p34`
- double `_p35`
- double `_p3_c4` [3]
 p_4 , 3-momentum of the two-photon central system
- double `_p3_l6` [3]
 p_6 , 3-momentum of the first outgoing lepton
- double `_p3_l7` [3]
 p_7 , 3-momentum of the second outgoing lepton
- double `_p3_p1` [3]
 p_1 , 3-momentum of the first proton-like incoming particle
- double `_p3_p2` [3]
 p_2 , 3-momentum of the second incoming particle

- double `_p3_p3` [3]
p₃, 3-momentum of the first proton-like outgoing particle
- double `_p3_p5` [3]
p₅, 3-momentum of the second proton-like outgoing particle
- double `_p45`
- double `_p_p3`
- double `_p_p4`
- double `_p_p5`
- double `_pc4`
|p₄|, 3-momentum norm of the two-photon central system
- int `_pdg1`
- int `_pdg2`
- int `_pdg3`
- int `_pdg5`
- int `_pdg6`
- int `_pdg7`
- double `_pl6`
|p₆|, 3-momentum norm of the first outgoing lepton
- double `_pl7`
|p₇|, 3-momentum norm of the second outgoing lepton
- double `_pp1`
|p₁|, 3-momentum norm of the first proton-like incoming particle
- double `_pp2`
|p₂|, 3-momentum norm of the second proton-like incoming particle
- double `_pp3`
|p₃|, 3-momentum norm of the first proton-like outgoing particle
- double `_pp5`
|p₅|, 3-momentum norm of the second proton-like outgoing particle
- double `_pt_l6`
p_{T,6}, transverse momentum of the first outgoing lepton
- double `_pt_l7`
p_{T,7}, transverse momentum of the second outgoing lepton
- double `_ptcut`
- double `_ptot`
Total momentum provided by the two incoming proton-like particles (along the z-axis)
- double `_q1dq`
- double `_q1dq2`
- double `_q2max`
Maximal Q² exchange.
- double `_q2min`
Minimal Q² exchange.
- double `_qp2max`
- double `_qp2min`
- double `_qve` [4]
- double `_s`
s, squared centre of mass energy of the incoming particles' system
- double `_s1`
- double `_s2`
- double `_sa1`
- double `_sa2`
- double `_sl1`
- double `_sp3`

- $\sin \phi_3$ of the first outgoing proton-like particle
- double [_sp5](#)
- $\sin \phi_5$ of the second outgoing proton-like particle
- double [_sp6](#)
- $\sin \phi_6$ of the first outgoing lepton
- double [_sp7](#)
- $\sin \phi_7$ of the second outgoing lepton
- double [_sqs](#)
- \sqrt{s} , centre of mass energy of the incoming particles' system
- double [_st3](#)
- $\sin \theta_3$ of the first outgoing proton-like particle
- double [_st4](#)
- $\sin \theta_4$ of the two-photons centre of mass system
- double [_st5](#)
- $\sin \theta_5$ of the second outgoing proton-like particle
- double [_st6](#)
- $\sin \theta_6$ of the first outgoing lepton
- double [_st7](#)
- $\sin \theta_7$ of the second outgoing lepton
- double [_stcm6](#)
- $\sin \theta_6^{\text{CM}}$, production angle of the first outgoing lepton, computed in the centre of mass system.
- double [_t1](#)
- double [_t2](#)
- double [_tau](#)
- double [_w1](#)
- m_1^2 , squared mass of the first proton-like incoming particle
- double [_w12](#)
- double [_w2](#)
- m_2^2 , squared mass of the second proton-like incoming particle
- double [_w3](#)
- m_3^2 , squared mass of the first proton-like outgoing particle
- double [_w31](#)
- double [_w4](#)
- m_4^2 , squared mass of the two-photon central system
- double [_w5](#)
- m_5^2 , squared mass of the second proton-like outgoing particle
- double [_w52](#)
- double [_w6](#)
- m_6^2 , squared mass of the first outgoing lepton
- double [_w7](#)
- m_7^2 , squared mass of the second outgoing lepton
- double [_wmax](#)
- double [_wmin](#)
- double * [_x](#)
- Array of `_ndim` components representing the point on which the weight in the cross-section is computed.
- bool [setin](#)
- Is the incoming particles' kinematic set ?
- bool [setkin](#)
- Is the full event's kinematic set ?
- bool [setll](#)
- Is the outgoing leptons' state set ?

- bool `setout`
Is the outgoing particles' kinematic set ?
- bool `setp1`
Is the first incoming proton-like particle's kinematic set ?
- bool `setp2`
Is the second incoming proton-like particle's kinematic set ?
- bool `setp3`
Is the first outgoing proton-like particle's kinematic set ?
- bool `setp5`
Is the second outgoing proton-like particle's kinematic set ?

2.2.1 Constructor & Destructor Documentation

2.2.1.1 GamGam::GamGam (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>x_[]</i>	The <i>ndim_</i> -dimensional point in the phase space on which the kinematics and the cross-section are computed

2.2.2 Member Function Documentation

2.2.2.1 void GamGam::ComputeSqS ()

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

Here is the caller graph for this function:



2.2.2.2 double GamGam::ComputeXsec (int *nm_* = 1)

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

Returns

$\sigma_{\gamma\gamma \rightarrow \ell^+\ell^-}$, the total cross section for the given event, provided the phase space region on which it is integrated.

2.2.2.3 bool GamGam::IsKinematicsDefined () [inline]

Is the system's kinematics well defined? Mandatory to perform the point's cross-section computation.

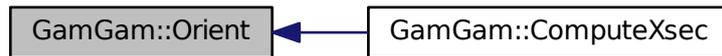
Returns

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

2.2.2.4 bool GamGam::Orient ()

Calculates energies and momenta of the 1st, 2nd, 3rd, 4th and 5th particle in the overall centre of mass frame.

Here is the caller graph for this function:

**2.2.2.5 double GamGam::PeriPP (int nup_, int ndown_)**

Contains the expression of the matrix element squared for the process under considerations. It returns the value of the product of the form factor or structure functions and the matrix element squared.

Returns

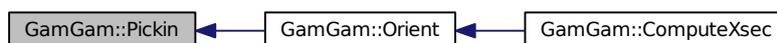
The full matrix element for the two-photon production of a pair of spin $-\frac{1}{2}$ -point particles

Here is the caller graph for this function:

**2.2.2.6 bool GamGam::Pickin ()**

Describes the kinematics of the process $p_1 + p_2 \rightarrow p_3 + p_4 + p_5$ in terms of Lorentz-invariant variables. Fills several variables to be used in PeriPP (essential for the evaluation of the matrix element).

Here is the caller graph for this function:



2.2.2.7 void GamGam::SetCuts (Cuts cuts_)

Parameters

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

2.2.2.8 bool GamGam::SetIncomingKinematics (int , double [], int)

Parameters

<i>part_</i>	Role of the particle in the process
<i>momentum_</i>	3-momentum of the particle
<i>pdgId_</i>	Particle ID according to the PDG convention

2.2.2.9 bool GamGam::SetOutgoingParticles (int *part_*, int *pdgId_*)

Parameters

<i>part_</i>	Role of the particle in the process
<i>pdgId_</i>	Particle ID according to the PDG convention

2.2.2.10 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.

2.2.3 Field Documentation

2.2.3.1 double GamGam::d1 [private]

[2] for the full definition of this quantity

2.2.3.2 double GamGam::d2 [private]

[2] for the full definition of this quantity

2.2.3.3 double GamGam::d4 [private]

[2] for the full definition of this quantity

2.2.3.4 double GamGam::d5 [private]

[2] for the full definition of this quantity

2.2.3.5 double GamGam::d6 [private]

[2] for the full definition of this quantity

2.2.3.6 double GamGam::d8 [private]

[2] for the full definition of this quantity

2.3 GnuPlot Class Reference

Public Member Functions

- **GnuPlot** (std::string outFile_="")

- int **DrawHistogram** ()
- int **Fill** (double)
- void **operator<<** (const std::string &command)
 - Feeds a command line to the [Gnuplot](#) interpreter.*
- void **SetHistogram** (int, double, double)
- void **SetLogy** (bool logy_=true)
 - Toggles the logarithmic scale for the y-axis.*
- void **SetOutputFile** (std::string)
- void **SetTitle** (std::string)
 - Sets the title for the graph.*
- void **SetXAxisTitle** (std::string)
 - Sets the caption for the x-axis.*
- void **SetYAxisTitle** (std::string)
 - Sets the caption for the y-axis.*

Protected Attributes

- FILE * **_pipe**
 - The pipe used to feed the [Gnuplot](#) interpreter.*

Private Attributes

- double * **__histBounds**
- double **__histHigh**
- double **__histLow**
- int **__histNum**
- int **__histOverflow**
- int **__histUnderflow**
- int * **__histValues**
- bool **__isHist**

2.3.1 Member Function Documentation

2.3.1.1 void Gnuplot::operator<< (const std::string & *command*)

Parameters

<i>&command</i>	The Gnuplot-formatted command line to feed
---------------------	--

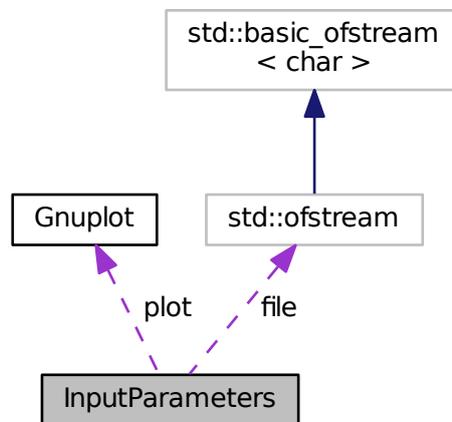
2.3.1.2 void Gnuplot::SetOutputFile (std::string *outFile_*)

Sets the file on which the graph has to be produced

2.4 InputParameters Class Reference

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

Collaboration diagram for InputParameters:



Data Fields

- bool **debug**
- 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 **maxpt**
Maximal transverse momentum of the outgoing leptons.
- int **mcut**
Set of cuts to apply on the outgoing leptons.
- double **minpt**
Minimal transverse momentum of the outgoing leptons.
- int **ncvg**
- int **p1mod**
First particle's mode.
- int **p2mod**
Second particle's mode.
- int **pair**
PDG id of the outgoing leptons.
- Gnuplot * **plot**

2.4.1 Field Documentation

2.4.1.1 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

2.4.1.2 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

2.4.1.3 int InputParameters::p2mod

Note

Was named EMOD in ILPAIR

2.4.1.4 int InputParameters::pair

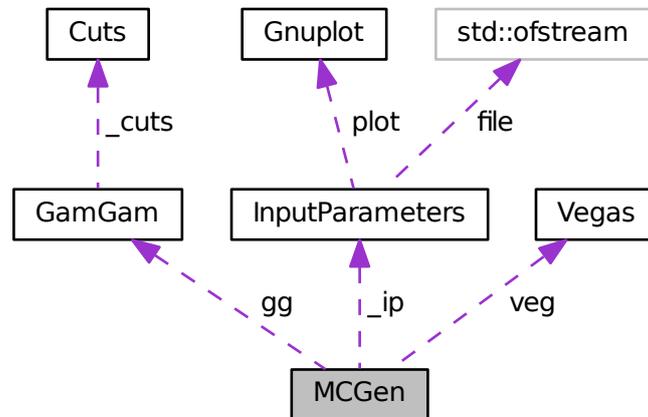
The particle code of produced leptons :

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

2.5 MCGen Class Reference

Core Monte-Carlo generator.

Collaboration diagram for MCGen:



Public Member Functions

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

Private Attributes

- int **_inp1pdg**
- int **_inp2pdg**
- [InputParameters _ip](#)
Set of parameters to setup the phase space to integrate.
- int **_ndim**
Number of dimensions on which to perform the integration.
- int **_outp1pdg**
- int **_outp2pdg**
- [GamGam * gg](#)
The GamGam object computing the kinematics and the cross-section for the given point in the phase space.
- [Vegas * veg](#)
The Vegas integrator which will integrate the function.

2.5.1 Constructor & Destructor Documentation

2.5.1.1 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

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

2.5.2 Member Function Documentation

2.5.2.1 InputParameters MCGen::GetInputParameters () [inline]

Returns

The InputParameter object embedded in this class

2.6 Particle Class Reference

Kinematics of one particle.

Public Member Functions

- `std::string GetLHEline ()`
- `void SetP (double, double, double)`
Sets the momentum.

Data Fields

- `double e`
Energy.
- `double m`
Mass.
- `int pdgId`
Particle Data Group integer identifier.
- `double pt`
Transverse momentum.
- `double px`
Momentum along the x-axis.
- `double py`
Momentum along the y-axis.
- `double pz`
Momentum along the z-axis.
- `int role`
Role in the considered process.

2.6.1 Detailed Description

Kinematic information for one particle

2.6.2 Member Function Documentation

2.6.2.1 `std::string Particle::GetLHEline ()`

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

Returns

The LHE line

2.6.2.2 void Particle::SetP (double px_, double py_, double pz_)

Parameters

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

2.7 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 `LaunchGeneration` (int)
Launches the generation of events.
- int `LaunchIntegration` ()
Launches the integration of the provided function.

Private Attributes

- `gsl_monte_function` * `_F`
The wrapped-up function to integrate, along with the input parameters.
- `size_t` `_ncalls`
Fixed number of function calls to use.
- `size_t` `_ndim`
The number of dimensions on which to integrate the function.
- `size_t` `_nlter`
Number of points to generate in order to integrate the function.
- `gsl_rng` * `_r`
GSL's random number generator.
- `gsl_monte_vegas_state` * `_s`
GSL's Vegas integration state structure.
- `double` * `_xl`
Lower bounds for the points to generate.
- `double` * `_xu`
Upper bounds for the points to generate.

2.7.1 Constructor & Destructor Documentation

2.7.1.1 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

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

2.7.2 Member Function Documentation

2.7.2.1 int Vegas::LaunchGeneration (int *nEvs_*)

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

Parameters

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

2.7.2.2 int Vegas::LaunchIntegration ()

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

References

- [1] G Peter Lepage. A new algorithm for adaptive multidimensional integration. *Journal of Computational Physics*, 27(2):192 – 203, 1978. [16](#)
- [2] J.A.M. Vermaseren. Two-photon processes at very high energies. *Nuclear Physics B*, 229(2):347 – 371, 1983. [10](#)

Index

- [_d1](#)
 - [GamGam, 10](#)
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 - [_d4](#)
 - [GamGam, 10](#)
 - [_d5](#)
 - [GamGam, 10](#)
 - [_d6](#)
 - [GamGam, 10](#)
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