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1: // Eikonal density and screening class
2: //
3: //
4: // (c) 2017-2019 Mikael Mieskolainen
5: // Licensed under the MIT License <http://opensource.org/licenses/MIT>.
6:
7: // C++
8: #include <fstream>
9: #include <future>
10: #include <iomanip>
11: #include <iostream>
12: #include <vector>
13: #include <cmath>
14:
15: // Own
16: #include "Graniitti/MAux.h"
17: #include "Graniitti/MEikonal.h"
18: #include "Graniitti/MForm.h"
19: #include "Graniitti/MMath.h"
20: #include "Graniitti/MTimer.h"
21: #include "Graniitti/MPDG.h"
22:
23: // Libraries
24: #include "json.hpp"
25: #include "rang.hpp"
26:
27:
28: using gra::aux::indices;
29: using gra::math::msqrt;
30: using gra::math::pow2;
31: using gra::math::zi;
32:
33:
34: namespace gra {
35:
36:     // Eikonal screening numerical integration parameters
37:     // N.B. Compile will make significant optimizations if boundaries
38:     // are const variables.
39:     namespace MEikonalNumerics {
40:
41:         constexpr double MinKT2 = 1E-6;
42:         constexpr double MaxKT2 = 25.0;
43:         unsigned int NumberKT2 = 0;
44:         bool logKT2 = false;
45:
46:         constexpr double MinBT = 1E-6;
47:         constexpr double MaxBT = 10.0 / PDG::GeV2fm;
48:         unsigned int NumberBT = 0;
49:         bool logBT = false;
50:
51:         constexpr double FBIntegralMinKT = 1E-9;
52:         constexpr double FBIntegralMaxKT = 30.0;
53:         constexpr unsigned int FBIntegralN = 10000;
54:
55:         constexpr double MinLoopKT = 1E-4;
56:         double MaxLoopKT = 1.75;
57:
58:         std::string GetHashString() {
59:             std::string str = std::to_string(MEikonalNumerics::MinKT2) +
60:                 std::to_string(MEikonalNumerics::MaxKT2) +
61:                 std::to_string(MEikonalNumerics::NumberKT2) +
62:                 std::to_string(MEikonalNumerics::logKT2) +
63:                 std::to_string(MEikonalNumerics::MinBT) +
64:                 std::to_string(MEikonalNumerics::MaxBT) +
65:                 std::to_string(MEikonalNumerics::NumberBT) +
66:                 std::to_string(MEikonalNumerics::logBT) +
67:                 std::to_string(MEikonalNumerics::FBIntegralMinKT) +
68:                 std::to_string(MEikonalNumerics::FBIntegralMaxKT) +
69:                 std::to_string(MEikonalNumerics::FBIntegralN)
70:             ;
71:             return str;
72:         }
73:
74:         void ReadParameters() {
75:
76:             // Read and parse
77:             using json = nlohmann::json;
78:
79:             const std::string inputfile = gra::aux::GetBasePath(2) + "/modeldata/" + "NUMERICS.json";
80:             const std::string data = gra::aux::GetInputData(inputfile);
81:             json j;
82:
83:             try {

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84:         j = json::parse(data);
85:
86:         // JSON block identifier
87:         const std::string XID = "NUMERICS_EIKONAL";
88:
89:         //MaxKT2 = j[XID]["MaxKT2"];
90:         NumberKT2 = j[XID]["NumberKT2"];
91:         logKT2 = j[XID]["logKT2"];
92:
93:         //MaxBT = j[XID]["MaxBT"]; MaxBT /= gra::math::GeV2fm; // Input as fermi, program
uses GeV^{-1}
94:         NumberBT = j[XID]["NumberBT"];
95:         logBT = j[XID]["logBT"];
96:
97:         //FBIntegralMaxKT = j[XID]["FBIntegralMaxKT"];
98:         //FBIntegralN = j[XID]["FBIntegralN"];
99:
100:         //MaxLoopKT = j[XID]["MaxLoopKT"];
101:
102:     } catch (...) {
103:         std::string str =
104:             "MEikonalNumerics::ReadParameters: Error parsing " + inputfile + " (Check for e
xtra/missing commas)";
105:         throw std::invalid_argument(str);
106:     }
107: }
108:
109: // -----
110:
111: // Screening loop (minimum values)
112: unsigned int NumberLoopKT = 15; // Number of kt steps
113: unsigned int NumberLoopPHI = 12; // Number of phi steps
114:
115: // User setup (ND can be negative, to get below the default)
116: void SetLoopDiscretization(int ND) {
117:     NumberLoopKT = std::max(3, 3 * ND + (int)NumberLoopKT);
118:     NumberLoopPHI = std::max(3, 3 * ND + (int)NumberLoopPHI);
119: }
120:
121: } // Namespace MEikonal ends
122:
123:
124: MEikonal::MEikonal() {
125: }
126:
127:
128: MEikonal::~MEikonal() {
129: }
130:
131:
132: // Return total, elastic, inelastic cross sections
133: void MEikonal::GetTotXS(double& tot, double& el, double& in) const {
134:     tot = sigma_tot;
135:     el = sigma_el;
136:     in = sigma_inel;
137: }
138:
139:
140: // Construct density and amplitude
141: void MEikonal::S3Constructor(double s_in, const std::vector<gra::MParticle>& initialstate_in, bool onlyeiko
nal) {
142:
143:     // This first
144:     MEikonalNumerics::ReadParameters();
145:
146:     // Mandelstam s and initial state
147:     s = s_in;
148:     INITIALSTATE = initialstate_in;
149:
150:     // -----
151:     // Calculate hash based on all free variables -> if something changed,
152:     // calculate new densities
153:
154:     // Proton density
155:     {
156:         std::cout << "Initializing <eikonal density> array:" << std::endl;
157:         MBT.sqrts = msqrt(s); // FIRST THIS
158:         MBT.Set("bt", MEikonalNumerics::MinBT, MEikonalNumerics::MaxBT, MEikonalNumerics::NumberBT,
MEikonalNumerics::logBT);
159:         MBT.InitArray(); // Initialize (call last!)
160:
161:         const std::string hstr = std::to_string(s) + std::to_string(INITIALSTATE[0].pdg) + "_" + st
d::to_string(INITIALSTATE[1].pdg) +

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162:                                     PARAM_SOFT::GetHashString() + MEikonalNumerics::GetHashString();
163:     const unsigned long hash = gra::aux::djb2hash(hstr);
164:     const std::string filename = gra::aux::GetBasePath(2) +
165:                                     "/eikonal/" + "MBT_" + std::to_stri
ng(INITIALSTATE[0].pdg) +
166:                                     "_" + std::to_string(INITIALSTATE[
1].pdg) +
167:                                     "_" + gra::aux::ToString(msqrt(s),0) + "_" + std::to_string(ha
sh);
168:
169:     // Try to read pre-calculated
170:     bool ok = MBT.ReadArray(filename);
171:
172:     while (!ok) { // Problem, re-calculate
173:         // Pointer to member function: ReturnType (ClassType::*)(ParameterTypes...)
174:         std::complex<double> (MEikonal::*f)(double) const = &MEikonal::S3Density;
175:         S3CalculateArray(MBT,f);
176:         MBT.WriteArray(filename, true);
177:         ok = MBT.ReadArray(filename);
178:     }
179:
180:
181:     // Calculate cross sections (is fast)
182:     S3CalcXS();
183:
184:     // Init cut Pomerons (is fast)
185:     S3InitCutPomerons();
186:
187:     // Amplitude
188:     if (onlyeikonal == false) {
189:         std::cout << "Initializing <eikonal amplitude> array:" << std::endl;
190:
191:         MSA.sqrts = msqrt(s); // FIRST THIS
192:         MSA.Set("kt2", MEikonalNumerics::MinKT2, MEikonalNumerics::MaxKT2, MEikonalNumerics::Number
KT2, MEikonalNumerics::logKT2);
193:         MSA.InitArray(); // Initialize (call last!)
194:
195:         const std::string hstr = std::to_string(s)
196:                                     + std::to_string(INITIALSTATE[0].pdg)
197:                                     + std::to_string(INITIALSTATE[1].pdg)
198:                                     + PARAM_SOFT::GetHashString()
199:                                     + MEikonalNumerics::GetHashString();
200:
201:         const unsigned long hash = gra::aux::djb2hash(hstr);
202:         const std::string filename = gra::aux::GetBasePath(2) + "/eikonal/" + "MSA_" +
203:                                     std::to_string(INITIALSTATE[0].pdg
) + "_" + std::to_string(INITIALSTATE[1].pdg) +
204:                                     "_" + gra::aux::ToString(msqrt(s),0) + "_" + std::to_string(ha
sh);
205:
206:     // Try to read pre-calculated
207:     bool ok = MSA.ReadArray(filename);
208:
209:     while (!ok) { // Problem, re-calculate
210:         // Pointer to member function: ReturnType (ClassType::*)(ParameterTypes...)
211:         std::complex<double> (MEikonal::*f)(double) const = &MEikonal::S3Screening;
212:         S3CalculateArray(MSA, f);
213:         MSA.WriteArray(filename, true);
214:         ok = MSA.ReadArray(filename);
215:     }
216:
217:     // Tag it done
218:     S3INIT = true;
219: }
220:
221:
222: // Proton bt-density by Fourier-Bessel transform of the t-density
223: // see <http://mathworld.wolfram.com/HankelTransform.html>
224: //
225: //
226: // For eikonalization see:
227: // [REFERENCE: Desgrolard, Giffon, Martynov, Predazzi, https://arxiv.org/abs/hep-ph/9907451v2]
228: // [REFERENCE: Desgrolard, Giffon, Martynov, Predazzi, https://arxiv.org/abs/hep-ph/0001149]
229: //
230: //
231: // For some discussion about Odderon versus Pomeron, see:
232: // [REFERENCE: Ewerz, Maniatis, Nachtmann, https://arxiv.org/abs/1309.3478]
233: std::complex<double> MEikonal::S3Density(double bt) const {
234:
235:     // Discretization of kt
236:     const double kt_STEP =
237:         (MEikonalNumerics::FBIntegralMaxKT - MEikonalNumerics::FBIntegralMinKT) /
238:         MEikonalNumerics::FBIntegralN;

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239:
240:     // N + 1!
241:     std::vector<std::complex<double>> f(MEikonalNumerics::FBIntegralN + 1, 0.0);
242:
243:     // Initial state configuration [NOT IMPLEMENTED = SAME FOR pp and ppbar]
244:     // pp
245:     if (INITIALSTATE[0].pdg == PDG::PDG_p && INITIALSTATE[1].pdg == PDG::PDG_p) {
246:         // TBD
247:     // ppbar
248:     } else if ((INITIALSTATE[0].pdg == PDG::PDG_p && INITIALSTATE[1].pdg == -PDG::PDG_p) ||
249:                (INITIALSTATE[0].pdg == -PDG::PDG_p && INITIALSTATE[1].pdg == PDG::PDG_p)) {
250:         // TBD
251:     }
252:
253:     // Loop over
254:     for (const auto& i : indices(f)) {
255:
256:         const double kt = MEikonalNumerics::FBIntegralMinKT + i * kt_STEP;
257:
258:         // negative, with Mandelstam  $t \sim -kt^2$ 
259:         const double t = -gra::math::pow2(kt);
260:
261:         // Proton form factors (could be here extended to multichannel)
262:         const double F_i = gra::form::S3F(t);
263:         const double F_k = gra::form::S3F(t);
264:
265:         // Pomeron trajectory  $\alpha(t)$ 
266:         const double alpha_P = gra::form::S3PomAlpha(t);
267:
268:         // Pomeron exchange amplitude:
269:         // [Regge signature x Proton Form Factor x coupling x Proton
270:         // Form Factor x coupling x Propagator ]
271:         const double s0 = 1.0; // Typical (normalization) scale  $\text{GeV}^{-2}$ 
272:
273:         // const std::complex<double> eta_0 =
274:         // std::exp(gra::math::zi*PARAM_SOFT::PHASE_O);
275:
276:         std::complex<double> A =
277:             gra::math::pow2(PARAM_SOFT::gN_P) * F_i * F_k *
278:             gra::form::ReggeEta(alpha_P, 1) *
279:             std::pow(s / s0, alpha_P - 1.0); // Pomeron (C-even)
280:
281:         // Value
282:         f[i] = A * gra::math::BESSJ0(bt * kt) * kt;
283:         // f[i] = A * std::cyl_bessel_j(0, bt * kt) * kt; // c++17
284:     }
285:     // Fourier-Bessel transformation denominator
286:     const double TD = 2.0*gra::math::PI;
287:
288:     return gra::math::CSIntegral(f, kt_STEP) / TD;
289: }
290:
291:
292: // Calculate elastic screening amplitude by "Eikonalization",
293: // in  $kt^2$  obtained via  $bt$ -space Fourier-Bessel integral
294: //
295: //
296: // Amplitude in  $bt$ -space:  $A_{el}(b_t) = i(1 - \exp(i*XI(b_t)/2))$ 
297: //
298: std::complex<double> MEikonal::S3Screening(double kt2) const {
299:
300:     // Local discretization
301:     const double STEP = (MEikonalNumerics::MaxBT - MEikonalNumerics::MinBT) /
302:                         MEikonalNumerics::FBIntegralN;
303:
304:     const double kt = gra::math::msqrt(kt2);
305:     std::vector<std::complex<double>> f(MEikonalNumerics::FBIntegralN + 1, 0.0);
306:
307:     // Numerical integral loop over impact parameter ( $b_t$ ) space
308:     for (const auto& i : indices(f)) {
309:
310:         const double bt = MEikonalNumerics::MinBT + i * STEP;
311:         const std::complex<double> XI = MBT.Interpolate1D(bt);
312:
313:         // I. STANDARD EIKONAL APPROXIMATION
314:         const std::complex<double> A =
315:             gra::math::zi * (1.0 - std::exp(gra::math::zi * XI / 2.0));
316:
317:         f[i] = A * gra::math::BESSJ0(bt * kt) * bt;
318:         // f[i] = A * std::cyl_bessel_j(0, bt * kt) * bt; // c++17
319:     }
320:     const double C = 2.0 * gra::math::PI; // phi-integral
321:

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322:         // Numerical integration
323:         return (2.0 * s) * C * gra::math::CSIntegral(f, STEP);
324:     }
325:
326:
327:     // Calculate screened total, elastic and inelastic cross sections
328:     // in the eikonal model
329:     //
330:     //
331:     //  $\int d^2b f(b) = \int_0^\infty \int_0^{2\pi} r dr d\theta f(r, \theta)$ 
332:     //  $= 2\pi \int_0^\infty f(r) r dr$ 
333:     //
334:     void MEikonal::S3CalcXS() {
335:         std::cout << "MEikonal::S3CalcXS:" << std::endl << std::endl;
336:
337:         // Local discretization
338:         const unsigned int N = 2 * 3000; // even number
339:         const double STEP = (MEikonalNumerics::MaxBT - MEikonalNumerics::MinBT) / N;
340:
341:         // Two channel eikonal eigenvalue solutions obtained via symbolic
342:         // calculation see e.g.
343:         //
344:         // [REFERENCE: Khoze, Martin, Ryskin, https://arxiv.org/abs/hep-ph/0007359v2]
345:         // [REFERENCE: Roehr, http://inspirehep.net/record/1351489/files/Thesis-2014-Roehr.pdf]
346:         //
347:         // Unitary transformation matrix
348:         const MMatrix<double> cc = {{ 1, 1, 1, 1}, // pp
349:                                     {-1, 1, 1, -1}, // pN*
350:                                     {-1, 1, -1, 1}, // N*p
351:                                     { 1, 1, -1, -1}}; // N*N*
352:
353:         // Eigenvalues
354:         const std::vector<double> lambda = {
355:             gra::math::pow2(1.0 - PARAM_SOFT::gamma),
356:             gra::math::pow2(1.0 + PARAM_SOFT::gamma),
357:             1.0 - gra::math::pow2(PARAM_SOFT::gamma),
358:             1.0 - gra::math::pow2(PARAM_SOFT::gamma)};
359:         // -----
360:         // N+1
361:         std::vector<std::complex<double>> f_tot(N + 1, 0.0);
362:         std::vector<std::complex<double>> f_el(N + 1, 0.0);
363:         std::vector<std::complex<double>> f_in(N + 1, 0.0);
364:
365:         // Two-channel eikonal, N+1!
366:         std::vector<std::vector<std::complex<double>>> f_2(
367:             4, std::vector<std::complex<double>>(N + 1, 0.0));
368:
369:         // Numerical integral loop over impact parameter (b_t) space
370:         for (const auto& i : indices(f_tot)) {
371:
372:             const double bt = MEikonalNumerics::MinBT + i * STEP;
373:
374:             // Calculate density
375:             const std::complex<double> XI = MBT.Interpolate1D(bt);
376:
377:             // -----
378:             // Single-Channel eikonal
379:             //
380:             // Elastic amplitude A_el(s,b)
381:             const std::complex<double> A_el = gra::math::zi * (1.0 - std::exp(gra::math::zi * XI / 2.0))
382:         };
383:
384:         // TOTAL: Im A_el(s,b)
385:         f_tot[i] = std::imag(A_el);
386:
387:         // ELASTIC: |A_el(s,b)|^2
388:         f_el[i] = gra::math::abs2(A_el);
389:
390:         // INELASTIC: 2Im A_el(s,b) - |A_el(s,b)|^2
391:         f_in[i] = 2.0 * std::imag(A_el) - gra::math::abs2(A_el);
392:         // -----
393:         // -----
394:         // Two-channel Eikonal solutions for pp->p(*)p(*)
395:         const std::vector<std::complex<double>> sol = {
396:             1.0 - std::exp(gra::math::zi * lambda[0] * XI / 2.0),
397:             1.0 - std::exp(gra::math::zi * lambda[1] * XI / 2.0),
398:             1.0 - std::exp(gra::math::zi * lambda[2] * XI / 2.0),
399:             1.0 - std::exp(gra::math::zi * lambda[3] * XI / 2.0)};
400:
401:         for (std::size_t k = 0; k < 4; ++k) {
402:
403:             // Amplitude squared

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404:         f_2[k][i] = gra::math::abs2(
405:             (sol[0] * cc[k][0] + sol[1] * cc[k][1] +
406:              sol[2] * cc[k][2] + sol[3] * cc[k][3]) / 4.0);
407:         f_2[k][i] *= bt; // Jacobian
408:     }
409:
410:     // Jacobian
411:     f_tot[i] *= bt;
412:     f_el[i] *= bt;
413:     f_in[i] *= bt;
414: }
415:
416: // 2D-Integral factor
417: const double IC = 2.0 * gra::math::PI;
418:
419: // Composite Simpson's rule, real is taken for C++ reasons in order to
420: // be able to substitute into double
421: sigma_tot = 2.0 * IC * std::real(gra::math::CSIntegral(f_tot, STEP)) * PDG::GeV2barn;
422: sigma_el = IC * std::real(gra::math::CSIntegral(f_el, STEP)) * PDG::GeV2barn;
423: sigma_inel = IC * std::real(gra::math::CSIntegral(f_in, STEP)) * PDG::GeV2barn;
424: // sigma_inel = sigma_tot - sigma_el; // cross check
425:
426: // Comments for the "multichannel" formalism [next version]
427:
428: // Total cross section: 2*\int d^2b \sum_{ik} |a_i|^2 |a_k|^2 (1 -
429: // exp^(-Omega(s,b)/2))
430: std::cout << "Single Channel Eikonal:" << std::endl;
431: printf(" Total xs:      %0.3f mb \n", sigma_tot*1E3);
432:
433: // Elastic cross section: \int d^2b (\sum_{ik} |a_i|^2 |a_k|^2 (1 -
434: // exp^(-Omega(s,b)/2))^2
435: printf(" Elastic xs:    %0.3f mb \n", sigma_el*1E3);
436:
437: // Inelastic cross section: \int d^2b \sum_{ik} |a_i|^2 |a_k|^2 (1 -
438: // exp^(-Omega(s,b)/2))
439: printf(" Inelastic xs:  %0.3f mb \n\n", sigma_inel*1E3);
440:
441: const double sigma_el_2 =
442:     IC * std::real(gra::math::CSIntegral(f_2[0], STEP)) * PDG::GeV2barn;
443: const double sigma_sd_a =
444:     IC * std::real(gra::math::CSIntegral(f_2[1], STEP)) * PDG::GeV2barn;
445: const double sigma_sd_b =
446:     IC * std::real(gra::math::CSIntegral(f_2[2], STEP)) * PDG::GeV2barn;
447: const double sigma_dd =
448:     IC * std::real(gra::math::CSIntegral(f_2[3], STEP)) * PDG::GeV2barn;
449:
450: // Scale
451: const double kappa = sigma_el / sigma_el_2;
452:
453: std::cout << "Two Channel normalized to Single Channel [PROTOTEST]" << std::endl;
454: printf("Calculated k = <el1>/<el2> = %0.3f \n", kappa);
455:
456: sigma_diff[0] = sigma_el_2 * kappa * 1E3;
457: sigma_diff[1] = sigma_sd_a * kappa * 1E3;
458: sigma_diff[2] = sigma_sd_b * kappa * 1E3;
459: sigma_diff[3] = sigma_dd * kappa * 1E3;
460:
461: printf(" pp   xs:  %0.3f mb \n", sigma_diff[0]);
462: printf(" pN*  xs:  %0.3f mb \n", sigma_diff[1]);
463: printf(" N*p   xs:  %0.3f mb \n", sigma_diff[2]);
464: printf(" N*N*  xs:  %0.3f mb \n\n", sigma_diff[3]);
465:
466: }
467:
468: // Calculate interpolation arrays
469: void MEikonal::S3CalculateArray(IArray1D& arr, std::complex<double> (MEikonal::*f)(double) const) {
470:     std::cout << "MEikonal::S3CalculateArray:" << std::endl;
471:     std::vector<std::future<std::complex<double>>> futures; // std::async return values
472:     MTimer timer(true);
473:
474:     // Loop over discretized variable
475:     for (std::size_t i = 0; i < arr.F.size_row(); ++i) {
476:
477:         const double a = arr.MIN + i * arr.STEP;
478:         arr.F[i][X] = a;
479:
480:         // Transform input to linear if log stepping, for the function
481:         const double var = (arr.islog) ? std::exp(a) : a;
482:
483:         gra::aux::PrintProgress(i / static_cast<double>(arr.N + 1));
484:         futures.push_back(std::async(std::launch::async, f, this, var));
485:     }
486:     gra::aux::ClearProgress();

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487:         std::cout << std::endl;
488:         printf("- Time elapsed: %0.1f sec \n\n", timer.ElapsedSec());
489:
490:         // Retrieve std::async values
491:         for (const auto& i : indices(futures)) {
492:             arr.F[i][Y] = futures[i].get();
493:         }
494:     }
495:
496: // Write the array to a file
497: bool IArray1D::WriteArray(const std::string& filename, bool overwrite) const {
498:
499:     // Do not write if file exists already
500:     if (gra::aux::FileExist(filename) && !overwrite) {
501:         // std::cout << "- Found pre-calculated" << std::endl;
502:         return true;
503:     }
504:     std::ofstream file;
505:     file.open(filename);
506:     if (!file.is_open()) {
507:         std::string str = "IArray1D::WriteArray: Fatal IO-error with: " + filename;
508:         throw std::invalid_argument(str);
509:     }
510:
511:     MTimer timer(true);
512:     std::cout << "IArray1D::WriteArray: ";
513:     for (std::size_t i = 0; i < F.size_row(); ++i) {
514:
515:         // Write to file
516:         file << std::setprecision(15)
517:             << std::real(F[i][X]) << ","
518:             << std::real(F[i][Y]) << ","
519:             << std::imag(F[i][Y]) << std::endl;
520:     }
521:     printf("Time elapsed %0.1f sec \n", timer.ElapsedSec());
522:     file.close();
523:     return true;
524: }
525:
526: // Read the array from a file
527: bool IArray1D::ReadArray(const std::string& filename) {
528:
529:     std::ifstream file;
530:     file.open(filename);
531:     if (!file.is_open()) {
532:         std::string str = "IArray1D::ReadArray: Fatal IO-error with: " + filename;
533:         return false;
534:     }
535:     std::string line;
536:     unsigned int fills = 0;
537:     std::cout << "IArray1D::ReadArray: ";
538:
539:     for (std::size_t i = 0; i < F.size_row(); ++i) {
540:
541:         // Read every line from the stream
542:         getline(file, line);
543:
544:         std::istringstream stream(line);
545:         std::vector<double> columns(3, 0.0);
546:         std::string element;
547:
548:         // Get every line element (3 of them) separated by separator
549:         int k = 0;
550:         while (getline(stream, element, ',')) {
551:             columns[k] = std::stod(element); // string to double
552:             ++k;
553:             ++fills;
554:         }
555:         F[i][X] = columns[0];
556:         F[i][Y] = std::complex<double>(columns[1], columns[2]);
557:     }
558:     file.close();
559:
560:     if (fills != 3 * (N + 1)) {
561:         std::string str = "Corrupted file: " + filename;
562:         std::cout << str << std::endl;
563:         return false;
564:     }
565:     std::cout << "[DONE]" << std::endl;
566:     return true;
567: }
568:
569:

```

```

570: // Standard 1D-linear interpolation
571: //
572: std::complex<double> IArray1D::Interpolate1D(double a) const {
573:
574:     const double EPS = 1e-5;
575:     if (a < MIN) { a = MIN; } // Truncate before (possible) logarithm
576:
577:     // Logarithmic stepping or not
578:     if (islog) { a = std::log(a); }
579:
580:     if (a > MAX*(1+EPS) ) {
581:         printf("IArray1D::Interpolate1D(%s) Input out of grid domain: "
582:             "%s = %0.3f [%0.3f, %0.3f] \n", name.c_str(), name.c_str(), a, MIN, MAX);
583:
584:         throw std::invalid_argument("Interpolate1D:: Out of grid domain");
585:     }
586:     int i = std::floor((a - MIN) / STEP);
587:
588:     // Boundary protection
589:     if (i < 0) { i = 0; } // Int needed for this, instead of unsigned int
590:     if (i >= (int)N) { i = N-1; } // We got N+1 elements in F
591:
592:     // y = y0 + (x - x0)*[(y1 - y0)/(x1 - x0)]
593:     return F[i][Y] + (a - F[i][X]) * ((F[i + 1][Y] - F[i][Y]) / (F[i + 1][X] - F[i][X]));
594: }
595:
596:
597: // Calculate the number of cut soft Pomerons for the inelastic
598: //
599: void MEikonal::S3InitCutPomerons() {
600:
601:     std::cout << "MEikonal::S3InitCutPomerons: [PROTOTEST]" << std::endl;
602:
603:     // Numerical integral loop over impact parameter (b_t) space
604:     const double STEP = (MEikonalNumerics::MaxBT - MEikonalNumerics::MinBT) / MEikonalNumerics::NumberB
T,
605:     P_array = std::vector<std::vector<double>>(MCUT, std::vector<double>(MEikonalNumerics::NumberBT+1,
0.0));
606:
607:     for (std::size_t j = 0; j < MEikonalNumerics::NumberBT+1; ++j) {
608:
609:         const double bt = MEikonalNumerics::MinBT + j*STEP;
610:         const double XI = std::imag(MBT.Interpolate1D(bt));
611:
612:         // Poisson probabilities P_m(bt)
613:         for (std::size_t m = 1; m < MCUT; ++m) {
614:
615:             // Poisson ansatz
616:             double P_m = std::pow(2*XI, m) / gra::math::factorial(m) * std::exp(-2*XI);
617:             P_array[m][j] = P_m * bt; // *bt from jacobian \int d^2b ...
618:         }
619:     }
620:
621:     // Impact parameter <bt> average probabilities
622:     P_cut.resize(MCUT, 0.0);
623:     for (std::size_t m = 0; m < MCUT; ++m) {
624:         P_cut[m] = gra::math::CSIntegral(P_array[m], STEP) / (MEikonalNumerics::MaxBT - MEikonalNum
erics::MinBT);
625:         printf("P_cut[m=%2lu] = %0.5f \n", m, P_cut[m]);
626:     }
627:     std::cout << "-----" << std::endl;
628:     printf("P_cut[SUM] = %0.5f \n", std::accumulate(P_cut.begin(), P_cut.end(), 0.0));
629:
630:     // Calculate zero-truncated average
631:     double avg = 0;
632:     for (std::size_t m = 1; m < P_cut.size(); ++m) {
633:         avg += m * P_cut[m];
634:     }
635:     printf("<P_cut[m>0> = %0.2f \n\n", avg);
636: }
637:
638: } // gra namespace ends
639:

```