"Once upon a time...

there was a Good Software Engineer whose Customers knew exactly what they wanted. The Good Software Engineer worked very hard to design the Perfect System that would solve all the Customers' problems now and for decades. When the Perfect System was designed, implemented and finally deployed, the Customers were very happy indeed. The Maintainer of the System had very little to do to keep the Perfect System up and running, and the Customers and the Maintainer lived happily every after."

It sounds like a fairy tale...

Could it be because there are no Good Software Engineers?

Could it be because the Customers/Users don't really know what they want?

Or because the Perfect System does not exist?

Lehman laws

M. M. Lehman, **Programs, Life Cycles, and Laws of Software Evolution,** *Proc. IEEE,* vol. 68, no. 9, Sep. 1980

1. Continuing Change

A program that is used and that as an implementation of its specification reflects some other reality, **undergoes continual change** or **becomes progressively less useful**. The change or decay process continues until it is judged more cost effective to replace the system with a recreated version.

2. Increasing Complexity

As an evolving program is continually changed, **its complexity**, *reflecting deteriorating structure*, **increases** unless work is done to maintain or reduce it.



Refactoring

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http://www.ge.infn.it/geant4/training/APC2025/



...is a disciplined technique for **improving the design of an existing code**

In the ideal world

there would be hardly any need for refactoring

In the real world of HEP (and related fields) most software needs to be refactored

By learning refactoring you also learn writing code that minimizes the need to be refactored

"If it ain't broken, don't fix it"

conventional wisdom

A piece of software can be broken in many ways



it no longer delivers the function it is designed to perform it can no longer be maintained

- Obsolete or no documentation
- Missing tests
- Original developers or users have left
- Inside knowledge about the system has disappeared
- Limited understanding of the entire system
- Too long to turn things over to **production**
- Too much time to make simple changes
- Need for constant bug fixes
- Big build times
- Difficulties separating products
- Duplicated code
- Code smells

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S. Demeyer, S. Ducasse, O. Nierstrasz, Object Oriented Reengineering Patterns

- Warnings you
 - are heading
 - into trouble
- usually do not occur isolated

Photoionisation cross section

Cross sections in Geant4 "standard" photoelectric model based on "improved" Biggs-Lighthill parameterisation

F. Biggs and R. Lighthill, Analytical Approximation for X-ray Cross Sections III, Sandia Lab. Report SAND-0070, 1988



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http://arxiv.org/abs/1601.06514





Post-RD44 electromagnetic design

Maria Grazia Pia, INFN Genova

Hidden dependencies

evolution



- 1. one of a set of prescribed movements
- 2. a process of change in a certain direction
- 3. the process of working out or developing
- 4. the historical development of a biological group
- 5. the extraction of a mathematical root
- 6. a process in which the whole universe is a progression of interrelated phenomena

"The code slowly sinks from engineering to hacking." M. Fowler, **Refactoring**



"With rapid development tools and rapid turnover in personnel, software systems can turn into legacies more quickly than you might imagine." S. Demeyer, S. Ducasse, O. Nierstrasz, Object Oriented Reengineering Patterns



- 1. a gift by will especially of money or other personal property
- something transmitted by or received from an ancestor or predecessor or from the past

"A legacy is something *valuable* that you have *inherited.*"

S. Demeyer, S. Ducasse, O. Nierstrasz, Object Oriented Reengineering Patterns

Refactoring

Reengineering

Methods and techniques

- to deal with software evolution
- to manage complexity
- to work with legacy code
- in a **disciplined** and **effective** way





Evolution of legacy systems



Take a loan on your software

 \Rightarrow pay back via reengineering

Investment for the future

 \Rightarrow paid back during maintenance

S. Demeyer, S. Ducasse, O. Nierstrasz, Object Oriented Reengineering Patterns

Software maintenance

"The modification of a software product after delivery to **correct faults**,

to **improve** performance or other attributes, or to **adapt** the product to a modified environment."

IEEE Standard 1219

- Accommodate changes in the software environment
- Incorporate new user requirements
- Fix errors
- Prevent future problems

OO techniques promise better	flexibility,
	reusability,
	maintainability
but they do not come for free!	•

Does refactoring pay back?

Steffen Hauf Refactoring Geant4 Radioactive Decay PhD Thesis

Well defined **responsibilities** and **interactions**





Motivations

- Gain understanding of the code
- Assess its capabilities and accuracy
- Improve physics performance
- Improve computational performance





Experiment: Z. W. Bell (ORNL)



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IEEE TRANSACTIONS ON NUCLEAR SCIENCE, VOL. 60, NO. 4, AUGUST 2013

Radioactive Decays in Geant4

Steffen Hauf, Markus Kuster, Matej Batič, Zane W. Bell, Dieter H. H. Hoffmann, Philipp M. Lang, Stephan Neff, Maria Grazia Pia, Georg Weidenspointner, and Andreas Zoglauer

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IEEE TRANSACTIONS ON NUCLEAR SCIENCE, VOL. 60, NO. 4, AUGUST 2013

Validation of Geant4-Based Radioactive Decay Simulation

Steffen Hauf, Markus Kuster, Matej Batič, Zane W. Bell, Dieter H. H. Hoffmann, Philipp M. Lang, Stephan Neff, Maria Grazia Pia, Georg Weidenspointner, and Andreas Zoglauer

Steffen Hauf Head of Software Control Group XFEL





Software technology

to deal with evolving/legacy software

- Problems
- Methods
- Techniques

Overview

Focus on basic concepts Guidance for further personal study



Peculiarities of refactoring physics software Hands-on practice

Common problems of legacy code

Insufficient documentation

- non-existent or out-of-date
- Improper layering
 - too few or too many layers
- Lack of modularity
 - strong coupling
- Duplicated code
 - copy & paste code

Duplicated functionality

similar functionality by separate teams

- **Misuse** of inheritance
 - code reuse vs. polymorphism
- Missing inheritance
 - duplication, case-statements
- Misplaced operations
 - operations outside classes
- Violation of encapsulation
 - type-casting; C++ "friends"
- Class abuse
 - classes as namespaces

S. Demeyer, S. Ducasse, O. Nierstrasz, Object Oriented Reengineering Patterns

1999





2018

Refactoring

is "the process of changing a software system in such a way that it does not alter the **external behavior** of the code yet improves its **internal structure**."

"When you refactor you are improving the design of the code after it has been written."

OBJECT-ORIENTED REENGINEERING PATTERN

Serge Demoper Stephane Ducasse Oscar Nierstrasz



Reengineering

Reengineering "seeks to transform a legacy system into the system you would have built if you had the luxury of hindsight and could have known all the new requirements that you know today."

"**Reengineering** [...] is the examination and alteration of a subject system to reconstitute it in a new form and the subsequent implementation of the new form."

"Reverse Engineering is the process of analyzing a subject system

- to identify the system's components and their interrelationships and
- create representations of the system in another form or at a higher level of abstraction."

"Forward Engineering is the traditional process of moving from highlevel abstractions and logical, implementation-independent designs to the physical implementation of a system."

E.J. Chikofsky, J.H. Cross, Reverse engineering and design recovery: a taxonomy, IEEE Software, vol. 7, no.1, pp. 13-17, 1990.

Basic concepts

Refactoring Reengineering

Interplay with testing

- What they are
- Why to refactor (reengineer)
- When to refactor
- What NOT to refactor
- When NOT to refactor

Do not refactor what does not pass the tests **Do not refactor** immediately before a critical deadline



To make software easier to understand and modify

When?

- Refactor when you want to add functionality
 - before adding it
- Refactor when you have to fix a bug
- Refactor while doing a code review (!)
- Refactor when you want to gain understanding of some legacy code
- **(**)...

Refactoring embedded in an iterative-incremental life-cycle

Set priorities while refactoring

Risks

- The new code may be more "elegant" but it has bugs / is slower
- Instead of just fixing a bug, you refactor the core code and screw up everything
 Testing
- When replacing old code that has been working fine for a long time, one risks reintroducing old problems that were fixed by some of the "ugly" (undocumented) code
- Nobody remembers all of the requirements, but break a single one by refactoring, and you can be in deep trouble
 Reengineering
- Refactoring increases the amount of verification testing that has to be done: when you refactor a class, you need to retest everything that deals with it (and side effects too)
 Planning
- Refactoring is an excuse for lazy programmers

Refactor your own code Refactor your colleague's code

What to refactor

How does one identify code that needs to be refactored?



Identifying code to be refactored



If it stinks, change it.

Grandma Beck, discussing child-rearing philosophy

M. Fowler, K. Beck et al., **Refactoring: Improving the Design of Existing Code**

A **code smell** is a **surface indication** that usually corresponds to a **deeper problem** in the system

Smells are heuristics that help in deciding:

- When to refactor
- What to refactor
- How to refactor

Quick to spot

Don't always indicate a problem



Code smells

- Duplicated code
- Long method
- Large class
- Long parameter list
- Divergent change
- Shotgun surgery
- Feature envy
- Data clumps
- Switch statement
- Parallel inheritance hierarchies
- Lazy class

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- Speculative generality
- Temporary field
- Comments
- Refused bequest
- Primitive obsession
- Message chains
- Middle man
- Inappropriate intimacy
- Alternative classes with different interfaces
- Incomplete library class
- Data class

M. Fowler, K. Beck et al.,

Refactoring: Improving the Design of Existing Code

Common code smells

Duplicated Code

- if you modify one instance of duplicated code but not the others, you may introduce a bug!

in the stink parade

Large class

- tries to do too much

Long Method

- difficult to understand and maintain
- Martin's Rule of Performance: Assume costs of lots of short functions are negligible and wait to be proven wrong!

Long Parameter List

- hard to understand, can become inconsistent

Code smells: dispensable

Data Class

- A data holder: a class that has attributes, getting and setting methods for the fields, and nothing else
- Objects should be about data and behavior

Speculative Generality

- "I may need the ability to do this kind of thing someday"

Lazy Class

- A class that no longer "pays its way" e.g. a class that was downsized by refactoring, or represented planned functionality that did not materialize

Dead Code

- Code that is not used

Code smells: OO abusers

Refused Bequest

 A subclass ignores most of the functionality provided by its superclass

Switch Statements

- Can be replaced by use of polymorphism

Temporary Field

- An attribute of an object is only set in certain circumstances *but an object should need all of its attributes*
- Fields used to hold intermediate results

Alternative Classes with Different Interfaces

- Two or more methods do the same thing but have different signature for what they do

Code smells: coupling

Middle Man

- A class delegates most of its responsibilities to another class
- *Does it really have a reason to exist?*

Message Chains

 A client asks an object for another object, then asks that object for another object etc.

Feature Envy

- A method requires lots of information from some other object

Inappropriate Intimacy

- Classes that know too much about each other's private details

Code smells: hindering change

Divergent Change

- Lack of cohesion:

one type of change requires changing one subset of methods; another type of change requires changing another subset

Shotgun Surgery

 A change requires lots of little changes in a lot of different objects

Parallel Inheritance Hierarchies

- Similar to Shotgun Surgery; each time I add a subclass to one hierarchy, I need to do it for all related hierarchies

...and more

Data Clumps

- Attributes that are used together, but are not part of the same object

Primitive Obsession

- A reluctance to use classes instead of primitive data types

Magic Number

- A literal value that appears in a program

Combinatorial Explosion

- Lots of code that does *almost* the same thing

How to refactor

Methods Techniques

Reverse engineering

Not limited to deriving a UML class diagram from the code... Motivation: understanding other people's code Refactoring is not meant to alter the behaviour of the code

To be tested!

Refactoring begins by designing a **solid set of tests** for the portion of code under analysis (usually unit tests)

Refactoring occurs as a series of small changes




Maria Graz

Geant4 version

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Then Endpoint Displacement of μ^{-} in the r ϕ Plane

geant4-10-00-ref-03, All MSC models, ARealisticRun, Gaussian fits



Maria Grazia Pia, INFN Genova



Fig. 8. Energy deposition divided by density-scaled calorimeter thickness Δx versus depth in C for 1 MeV e- at 0° incidence: simulations with Geant4 library-based processes, Geant4 version 8.1p02 (triangles) and 9.1 (circles), experimental data from [5] (crosses).



IEEE NSS Best Student Paper, 2007

IEEE TRANSACTIONS ON NUCLEAR SCIENCE, VOL. 60, NO. 4, AUGUST 2013

Validation of Geant4 Simulation of Electron **Energy Deposition**

Matej Batič, Gabriela Hoff, Maria Grazia Pia, Paolo Saracco, and Georg Weidenspointner



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Refactoring aims to preserve correctness

Sweeping under the carpet?



Was the original code verified?

Was the original code validated?



IEEE Standard 1012 Software Verification & Validation

What was the test coverage?

Were the test process and the test results documented?

Std 1012 evolution IEEE Standard for <u>System</u>, Software, and <u>Hardware</u> Verification and Validation

Risk mitigation

Do not mix!

One of the motivations for refactoring may be the need to introduce new features in the code

Refactoring

Does not modify the code behaviour



Adding new features

Modifying the code behaviour

- Test
- Refactor
- Test
- Add new feature
- Test
- Add new feature
- Test



Basic actions for common smells

Method invocation

- Consolidate recurring code into a single method
- OK when the recurring code does not span methods and all methods containing code belong to the same class



Inheritance

- Common code in two different classes
- New superclass introduced

class Child : public Super {
...
};

Extract commonality

Abstract interfaces are classes with no implementation Abstract classes

represent mixed design and implementation

Introduce abstract class

 a class with no or partial implementation

Add delegation

 Delegate the recurring code segments to a helper class class Common {
 void commonCode(...) {...}
 virtual void contextSpecificCode () = 0;
 ...
};

class Extension : public Common {
 void method1(...) {

helper.SomeMoreCommonCode();

}

...sounds like common sense?

Many refactoring techniques are just good practices of code hygiene

Apply them when writing new code!

The "legacy code" you may have to refactor in a few months/years may **be your own**... A colleague collaborating at your project may have to refactor your "legacy code"...

Refactoring Techniques

Composing Methods

Extract Method, Inline Method,...

Moving Features Between Objects Move Method, Move Field, Hide Delegate, ...

Organizing Data

Replace Data Value with Object, ...

Simplifying Conditional Expressions

Decompose Conditionals, ...

http://www.refactoring.com/catalog/index.html







- Same expression in two methods of the **same class**
 - Use **Extract Method** refactoring
- Same expression in two methods of sibling classes
 - Use Extract Method and Pull Up Method
- If code is similar, but not same
 - Consider Form Template Method
- Duplicated code in unrelated classes
 - May need to **Extract Class**
 - Or eliminate one of the versions

G4WentzelVIModel.cc

G4WentzelVIRelModel.cc

G4double G4WentzelVIModel::ComputeGeomPathLength(G4double truelength)	G4double G4WentzelVIRelMo	<pre>del::ComputeGeomPathLength(G4double truelength)</pre>
<pre>{ tPathLength = truelength; zPathLength = tPathLength;</pre>	{	th; gth;
<pre>if(lambdaeff > 0.0 && lambdaeff != DBL_MAX) { G4double tau = tPathLength/lambdaeff; //G4cout << "ComputeGeomPathLength: tLength= " << tPathLength // << " Leff= " << lambdaeff << " tau= " << tau << G4endl; // small step if(tau < numlimit) { zPathLength *= (1.0 - 0.5*tau + tau*tau/6.0); } }</pre>	<pre>if(lambdaeff > 0.0 && l G4double tau = tPathL //G4cout << "ComputeG // << "Leff= " << // small step if(tau < numlimit) { zPathLength *= (1.0 } }</pre>	<pre>ambdaeff != DBL_MAX) { ength/lambdaeff; eomPathLength: tLength= " << tPathLength lambdaeff << " tau= " << tau << G4endl;</pre>
<pre>// medium step } else { G4double e1 = 0.0; if(currentRange > tPathLength) { e1 = GetEnergy(particle,currentRange-tPathLength,currentCouple); } e1 = 0.5*(e1 + preKinEnergy); cosTetMaxNuc = wokvi->SetupKinematic(e1, currentMaterial); lambdaeff = GetTransportMeanFreePath(particle,e1); zPathLength = lambdaeff*(1.0 - G4Exp(-tPathLength/lambdaeff)); } else { lambdaeff = DBL_MAX; } //G4cout<<"Comp.geom: zLength= "<<zpathlength< td=""></zpathlength<></pre>	<pre>// medium step } else { G4double e1 = 0.0; if(currentRange > t e1 = GetEnergy(partic } e1 = 0.5*(e1 + preK cosTetMaxNuc = wokv lambdaeff = GetTran zPathLength = lambd } } else { lambdaeff = DB //G4cout<<"Comp.geom: z return zPathLength; }</pre>	<pre>PathLength) { le,currentRange-tPathLength,currentCouple); inEnergy); i->SetupKinematic(e1, currentMaterial); sportMeanFreePath(particle,e1); aeff*(1.0 - G4Exp(-tPathLength/lambdaeff)); L_MAX; } Length= "<<zpathlength<<" "<<tpathlength<<g4endl;<="" pre="" tlength=""></zpathlength<<"></pre>

Identical member function!

G4PAIModel.cc

G4PAIPhotModel.cc

G4PAIModel.cc vs. G4PAIPhotModel.cc								
G4PAIModel.cc - /Users/pia/Documents/g4dev/g4_svn/geant4.10.00.p01/source/processes/electromagnetic/standard/src G4PAIPhotModel.cc - /Users/pia/Documents/g4dev/g4_svn/geant4.10.00.p01/source/processes/electromagnetic/standard/src								
G4PAIModel::G4PAIModel()	G4PAIPhotModel::G4PAIPhotModel()							
<u>+</u>								
void G4PAIModel::Initialise(const G4ParticleDefinition* p.	11 void G4PAIPhotModel::Initialise(const G4ParticleDefinition* p, const G4DataVector& cuts)							
const G4DataVector& cuts)	{ { // if(f)/arbase > 0)							
if(fVerbose > 0) {	12 { Caputo all CAPATINe the data is the track in the second sec							
G4COUT<<"G4PAIMODEL::Initialise for "< <p->GetParticleName()<<g4enol; }</g4enol; </p->	G4COUt<<"G4PAIPhotmode(::Initialise for "< <p->GetParticleName()<<g4end(; }</g4end(; </p->							
if(isInitialised) { return; } isInitialised = true;	<pre>if(isInitialised) { return; } isInitialised = true;</pre>							
<pre>SetParticle(p); fParticleChange = GetParticleChangeForLoss();</pre>	<pre>SetParticle(p); fParticleChange = GetParticleChangeForLoss();</pre>							
<pre>if(IsMaster()) {</pre>	<pre>if(IsMaster()) {</pre>							
<pre>InitialiseElementSelectors(p, cuts);</pre>	<pre>InitialiseElementSelectors(p, cuts);</pre>							
if(!fModelData) {	if(!fModelData) {							
G4double tmin = LowEnergyLimit()∗fRatio; G4double tmax = HighEnergyLimit()∗fRatio; fModelData = new G4PAIModelData(tmin, tmax, fVerbose);	G4double tmin = LowEnergyLimit()*fRatio; G4double tmax = HighEnergyLimit()*fRatio; 13 ModelData = new G4PAIPhotData(tmin, tmax, fVerbose);							
<pre>} // Prepare initialization const G4MaterialTable≠ theMaterialTable = G4Material::GetMaterialTable(); size_t numOfMat = G4Material::GetNumberOfMaterials(); size_t numRegions = fPAIRegionVector.size();</pre>	<pre>} // Prepare initialization const G4MaterialTable* theMaterialTable = G4Material::GetMaterialTable(); size_t numOfMat = G4Material::GetNumberOfMaterials(); size_t numRegions = fPAIRegionVector.size();</pre>							
<pre>for(size_t iReg = 0; iReg < numRegions; ++iReg) { const G4Region* curReg = fPAIRegionVector[iReg]; G4Region* reg = const_cast<g4region*>(curReg);</g4region*></pre>	<pre>for(size_t iReg = 0; iReg < numRegions; ++iReg) { const G4Region* curReg = fPAIRegionVector[iReg]; G4Region* reg = const_cast<g4region*>(curReg);</g4region*></pre>							
<pre>for(size_t jMat = 0; jMat < numOfMat; ++jMat) { G4Material& mat = (*theMaterialTable)[jMat]; const G4MaterialCutsCouple& cutCouple = reg->FindCouple(mat); //G4cout << "Couple <" <<futcouple <<="" pre=""> (</futcouple></pre>	<pre>for(size_t jMat = 0; jMat < numOfMat; ++jMat) { G4Material* mat = (*theMaterialTable)[jMat]; const G4MaterialCutsCouple* cutCouple = reg->FindCouple(mat); //G4cout << "Couple <" << fCutCouple <" (set and the formation of t</pre>							
<pre>// GCOUT << CUTS.SIZE() << GAENOL; */ // check if this couple is not already initialized size_t n = fMaterialCutsCoupleVector.size(); if(0 < n) { for(size_t i=0; i<fmaterialcutscouplevector.size(); ++i)="" break;="" if(cutcouple="fMaterialCutsCoupleVector[i])" pre="" {="" }="" }<=""></fmaterialcutscouplevector.size();></pre>	<pre>// 64Cout << cuts.size() << 64enut; */ // check if this couple is not already initialized size_t n = fMaterialCutsCoupleVector.size(); if(0 < n) { for(size_t i=0; i<fmaterialcutscouplevector.size(); ++i)="" <="" break;="" if(cutcouple="fMaterialCutsCoupleVector[i])" pre="" {="" }=""></fmaterialcutscouplevector.size();></pre>							
<pre>} } // initialise data banks // initialise data banks fMaterialCutsCoupleVector.push_back(cutCouple); G4double deltaCutInKinEnergy = cuts[cutCouple->GetIndex()]; fModelData->Initialise(cutCouple, deltaCutInKinEnergy, this); }</pre>	<pre>} } // initialise data banks fMaterialCutsCoupleVector.push_back(cutCouple); G4double deltaCutInKinEnergy = cuts[cutCouple->GetIndex()]; fModelData->Initialise(cutCouple, deltaCutInKinEnergy, this); }.</pre>							
status: 44 differences IGENTICAI DUT THE INSTANTIO	ation of a "model data" object 🗸							

How to find duplicated code?

- Automated tools
 - Some exist
- By hand
 - Still the most common way
 - Not necessarily the most efficient (nor the most inefficient)

Reverse engineering

- Gain understanding of the code

G4KleinNishinaCompton.cc - /Users/pia/Documents/g4dev/g4_svn/geant4.10.00.p01/source/processes/electromagnetic/standard/src

G4KleinNishinaCompton::G4KleinNishinaCompton()

G4KleinNishinaCompton.cc

G4KleinNishinaModel.cc - /Users/pia/Documents/g4dev/g4_svn/geant4.10.00.p01/source/pro

G4KleinNishinaModel::G4KleinNishinaModel()

G4KleinNishinaModel.cc

Zddouble CAKLEINNIShinaCompton::Computed:ressSectionPerAted ///ooo000000000oo000000000	//		, ,
Gdouble Gmetherery, Gdouble Z, Gdouble,Gdouble Section = 0.0; If (Z < 0.9090)	G4double G4KleinNishinaCompton::ComputeCrossSectionPerAtom(//000000000000000000000
Gddouble GammaEnergy, Gddouble, Call, Gddouble, Gddoubl	const G4ParticleDefinition*,		
Gddouble Z, Gddouble, Gddouble Z, Gddouble, Gddouble Scction = 0.0; If (Z < 0.999) / return Section; If (Z < 0.999) / return Section; Static const Gddouble section = 0.0; If (Z < 0.999) / return Section; Static const Gddouble x = 20.0, b = 230.0, c = 440.0; Gddouble piZ = 2x(d1 + eixZ + fizXZ), piZ = 2x(d2 + eizZ + fizXZ); piZ = 2x(d1 + eizZ + fizXZ), piZ = 2x(d2 + eizZ + fizXZ); piZ = 2x(d1 + eizZ + fizXZ), piZ = 2x(d2 + eizZ + fizXZ); Gddouble piZ = 2x(d1 + eizZ + fizXZ), piZ = 2x(d2 + eizZ + fizXZ); piZ = 2x(d1 + eizZ + fizXZ), piZ = 2x(d2 + eizZ + fizXZ); Gddouble r0 = 15.04keV; If (Z < 1.5) T0 = 40,04keV; If (Z < 1.5) T0 = 40,04ke	G4double GammaEnergy,	17	G4double
Gddouble, sedouble) Gddouble, sedouble) Gddouble xSection = 0.0; if (Z < 0.9999)	G4double Z, G4double,	>	G4KleinNishinaModel::ComputeCrossSectionPerAtom(const G4ParticleDefinition*,
Gdouble section = 0.0; If (2 < 0.599)	G4aouble, G4aouble)		G4double GammaEnergy,
<pre>If (2 < 0.9999)</pre>	64 double xSection = 0.0 :		G4double, G4double)
if (GammaEnergy < 0.1+keV) / return xSection;	if (Z < 0.9999) return xSection;		{
<pre>// if (GamaEnergy > (100.sGeV/Z)) return xSection; static const G4double a = 20.0, b = 230.0, c = 440.0; G4double pIZ = Z*(d1 + e1z + f1+Z*Z), pZZ = Z*(d2 + e2+Z + f2*Z*Z); pJZ = Z*(d2 + e2+Z + f1*Z*Z), pZZ = Z*(d2 + e2+Z + f2*Z*Z); G4double T0 = 15.0*keV; if (Z < 1.5) T0 = 40.0*keV; G4double T0 = 15.0*keV; if (Z < 1.5) T0 = 40.0*keV; if (GamaEnergy, T0) / electron_mass_c2; xSection = p1Z*64Log(1.+2,*X)/X + (p2Z + p3Z*x + p4Z*xeX)/(1. + a*X + b*X*X + c*X*xX); // modification for low energy. (special case for Hydrogen) if (GamaEnergy < T0) { G4double f0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; G4double f0 = 1.*keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = 1.*keV; x = (T0+dT0) / electron_mass_c2; G4double f0 = 1.*keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f0 = keV; if (Z > 1.5) (c 2 = 0.15); d4double f</pre>	if (GammaEnergy < 0.1*keV) return xSection;	18	G4double xSection = 0.0 ;
<pre>static const G4double a = 20.0, b = 230.0, c = 440.0; G4double p12 = 2x(d1 + e1x + 1;1x2x2), p22 = 2x(d2 + e2x2 + 1;2x2x2), p32 = 2x(d3 + e3x2 + 1;3x2x2), p42 = 2x(d4 + e4x2 + 1;4x2x2); G4double p1 = 15.0+kev; if (2 < 1.5) T0 = 40.0+kev; if (2 < 1.5</pre>	// if (GammaEnergy > (100.*GeV/Z)) return xSection;		if (Z < 0.9999 GammaEnergy < 0.1*keV) { return xSection; }
Static const Gudouble a 2 0.0, b = 20.0, b =			
G4double plz = 2*(d1 + e1*2 + f1*2*2), plz = 2*(d2 + e2*2 + f2*2*2), plz = 2*(d3 + e3*2 + f3*2*2), pdz = 2*(d4 + e4*2 + f4*2*2); G4double plz = 2*(d1 + e1*2 + f1*2*2), pdz = 2*(d4 + e4*2 + f4*2*2); G4double plz = 15.9*keV; if (2 < 1.5) T0 = 40.0*keV; G4double X = max(GammaEnergy, T0) / electron_mass_c2; xSection = plz*G4Log(1.+2.*X)/X + (p22 + p32*X + p42*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double dT0 = 1.*keV; x = (T0+10) / electron_mass_c2; G4double c1 = -T0*(sigma-xSection)/(xSection*dT0); G4double c2 = 0.150; if (2 > 1.5) c2 = 0.375-0.0555*64Log(2); G4double c2 = 0.150; if (2 > 1.5) c2 = 0.375-0.0555*64Log(2); G4double y = G4Log(GamaEnergy, T0); xSection *= G4Exp(-y*(cl+c2*y)); }// G4cout < "e= " < GammaEnergy < "Z = " < Z <" cross= " < xSection < G4end1; return xSection; //ooo000000000oo000000000	static const G400uble $a = 20.0$, $b = 230.0$, $c = 440.0$;	19	static const 64double a = 20.0 , b = 230.0 , c = 440.0;
<pre>Solution p3z = Zx(d3 + e3xZ + f3xZxZ), p4Z = Zx(d4 + e4xZ + f4xZxZ); G4double T0 = 15,0*keV; if (Z < 1,5) f 0 = 40.0*keV; if (Z < 1,5) [T0 = 40.0*keV;</pre>	G4dauble n17 = 7*(d1 + e1*7 + f1*7*7), n27 = 7*(d2 + e2*7 + f2*7*7).		64double p17 = 7*(d1 + e1*7 + f1*7*7), p27 = 7*(d2 + e2*7 + f2*7*7).
G4double T0 = 15.0*keV; if (Z < 1.5) T0 = 40.0*keV;	p3Z = Z*(d3 + e3*Z + f3*Z*Z), p4Z = Z*(d4 + e4*Z + f4*Z*Z);		$p_{3Z} = Z*(d3 + e_{3*Z} + f_{3*Z*Z}), p_{4Z} = Z*(d4 + e_{4*Z} + f_{4*Z*Z});$
G4double T0 = 15.0*keV; if (Z < 1.5) T0 = 40.0*keV;			
<pre>if (Z < 1.5) T0 = 40.0*keV; G4double X = max(GammaEnergy, T0) / electron_mass_c2; xSection = p12*64Log(1+2,*x)/X + (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X); // modification for low energy. (special case for Hydrogen) if (GammaEnergy < T0) { G4double dT0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; (G4double dT0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; (G4double dT0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; (G4double dT0 = keV; X = (T0+dT0) / electron_mass_c2; (G4double dT0 = keV; (T = -T0*elsigna-xSection)/(xSection*dT0); (G4double dT0 = keV; (T</pre>	G4double T0 = 15.0*keV;	20	G4double T0 = 15.0*keV;
64double X = max(GammaEnergy, T0) / electron_mass_c2; xSection = p12+64Log(1,+2,+X)/X + (p22 + p32*X + p42*XX)/(1, + a*X + b*X*X + c*X*XX); // modification for low energy. (special case for Hydrogen) if (GammaEnergy < T0) {	if (Z < 1.5) TØ = 40.0*keV;		if (Z < 1.5) { T0 = 40.0*keV; }
<pre>Verdeduct x = max(cdum_mas_ct; xSection = p12x6dLog(1.+2.*x)/X + (p2Z + p32xX + p42*X*X)/(1. + a*X + b*X*X + c*X*X*X); // modification for low energy. (special case for Hydrogen) if (GammaEnergy < T0 { G4double dT0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; G4double sigma = p12x6dLog(1.+2*X)/X + (p2Z + p32*X + p42*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double c1 = -T0*(sigma-XSection)/(XSection*dT0); G4double c2 = 0.150; if (Z > 1.5) c2 = 0.375-0.9556*G4Log(Z); G4double y = G4Log(GammaEnergy/T0); XSection * G4Exp(-y*(c1+c2*y)); } // G4double y = G4Log(GammaEnergy/T0); XSection * G4Exp(-y*(c1+c2*y)); } // G4double y = G4Log(GammaEnergy/T0); XSection * G4Exp(-y*(c1+c2*y)); } // G4double y = G4Log(GammaEnergy < " Z= " << Z <" cross= " << xSection << G4endl; return xSection; } //ooo000000000ooo000000000ooo000000</pre>	6/double X = max(GammaEpergy T0) / electron mass co		(Adouble X - max(GammaEpergy TA) / electron mass c2:
<pre>+ (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X); // modification for low energy. (special case for Hydrogen) if (GammaEnergy < T0) { G4double dT0 = 1.**key; X = (T0+dT0) / electron_mass_c2; G4double dT0 = 1.**key; + (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double dT0 = 1.**key; x = (T0+dT0) / electron_mass_c2; G4double dT0 = 1.**key; + (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double c1 = -T0*(sigma-Xsection)/(Xsection*dT0); G4double c2 = 0.150; if (Z > 1.5) c2 = 0.375-0.0556*G4Log(Z); G4double c2 = 0.375-0.0556*G4Log(Z); if (Z > 1.5) { c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); G4couble c2 = 0.375-0.0556*G4Log(Z); if (X > 1.5) { c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); if (X > 1.5) { c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); if (X > 1.5) { c2 = 0.375-0.0556*G4Log(Z); } G4couble c2 = 0.375-0.0556*G4Log(Z); } C4couble c2 = 0.375-0.0556*</pre>	x section = p17×641 or (1+2, *X)/X		x Section = $p17 \times 64 \log(1 + 2 \times X)/X$
<pre>// modification for low energy. (special case for Hydrogen) if (GammaEnergy < T0) { G4double dT0 = 1.*keV; X = (Tpe4T0) / electron_mass_c2; G4double sigma = p12*G4Log(1.+2*X)/X</pre>	+ (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X);		+ $(pZZ + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X);$
<pre>// modification for low energy. (special case for Hydrogen) if (GammaEnergy < TO) { G4double dT0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; G4double sigma = p12*64Log(1.+2*X)/X + (p22 + p32*X + p42*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double c2 = 0.150; if (Z > 1.5) c2 = 0.375-0.0556*G4Log(2); G4double y = 64Log(CammaEnergy <<" Z= " << Z <<" cross= " << xSection *= G4Exp(-y*(c1+c2*y)); } // G4cout <<"e=" << GammaEnergy <" Z= " << Z <<" cross= " << xSection = 0.0; // G4cout <<"e=" << GammaEnergy <" Z= " << Z //ooo0000000000ooo0000000000</pre>			
<pre>11 (GammaEnergy < 10) { G4double dT0 = 1.*keV; X = (T0+dT0) / electron_mass_c2; G4double sigma = p1Z*64Log(1.+2*X)/X</pre>	// modification for low energy. (special case for Hydrogen)		<pre>// modification for low energy. (special case for Hydrogen)</pre>
<pre>tdouble did = 1*key; X = (T0+dT0) / electron_mass_c2; G4double sigma = p12k64Log(1.+2*X)/X</pre>	11 (GammaEnergy < 10) {	21	1† (GammaEnergy < T0) {
<pre>X = (10+010) / etection_mass_12, 2, G4double sigma = p12x64Log(1,+2*X)/X + (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double c1 = -T0*(sigma-xSection)/(xSection*dT0); G4double c2 = 0.150; if (Z > 1.5) c2 = 0.375-0.0556*G4Log(Z); G4double y = G4Log(GammaEnergy/T0); xSection *= G4Exp(-y*(c1+c2*y)); } // G4cout << "e= " << GammaEnergy << " Z= " << Z << " cross= " << xSection << G4end1; return xSection; } //ooo000000000ooo000000000ooo000000</pre>	G400DDLC 010 = 1.*KCV;		V = (T0+dT0) / electron mass c2 :
<pre>+ (p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X); G4double c1 = -T0*(sigma-xSection)/(xSection*dT0); G4double c2 = 0.150; if (Z > 1.5) c2 = 0.375-0.0556*G4Log(Z); G4double y = G4Log(GammaEnergy/T0); xSection *= G4Exp(-y*(c1+c2*y)); } // G4cout << "e= " << GammaEnergy <<" Z= " << Z <<" cross= " << xSection << G4endl; return xSection; } //ooo000000000ooo000000000ooo000000</pre>	$A = (1+\alpha) = 17 + 64 \ln \alpha (1+2*)/X$		$(10+010)$ / electron_mass_c2 , G4double sigma = p17*G4Log(1.+2*X)/X
G4double c1 = -T0*(sigma-xSection)/(xSection*dT0); G4double c2 = 0.150; if (z > 1.5) c2 = 0.375-0.0556*G4Log(Z); G4double y = G4Log(GammaEnergy/T0); xSection *= G4Exp(-y*(c1+c2*y)); } // G4cout <	+ $(p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X);$		+ $(p2Z + p3Z*X + p4Z*X*X)/(1. + a*X + b*X*X + c*X*X*X);$
G4double c2 = 0.150; G4double c2 = 0.150; if (Z > 1.5) c2 = 0.375-0.0556*G4Log(Z); G4double y = G4Log(GammaEnergy/T0); xSection *= G4Exp(-y*(c1+c2*y)); Section *= G4Exp(-y*(c1+c2*y)); // G4cout << "e= " << GammaEnergy << " Z= " << Z << " cross= " << xSection << G4endl;	G4double c1 = -T0*(sigma-xSection)/(xSection*dT0);		G4double c1 = -T0*(sigma-xSection)/(xSection*dT0);
<pre>if (Z > 1.5) {C2 = 0.375-0.0556*64Log(Z); G4double y = G4Log(GammaEnergy/T0); xSection *= G4Exp(-y*(c1+c2*y)); } // G4cout << "e= " << GammaEnergy << " Z= " << Z << " cross= " << xSection << G4endl; return xSection; } //ooo000000000ooo000000000 } // G4cout << "e= " << GammaEnergy << " Z= " << Z <!-- // G4cout << "e= " << GammaEnergy << " Z= " << Z // // G4cout << "e= " << GammaEnergy << " Z= " << Z // // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout << " e= " << GammaEnergy << " Z= " << Z // G4cout <</ re--></pre>	G4double c2 = 0.150;	22	G4double c2 = 0.150;
Gdouble y = G4Log(GammaEnergy/10); xSection *= G4Exp(-y*(c1+c2*y)); // G4cout <	if (Z > 1.5) c2 = 0.375-0.0556*64Log(Z);		if $(Z > 1.5)$ { $c2 = 0.375 - 0.0556 * G4Log(Z)$; }
<pre>xsection *= 64Exp(-y*(C1+C2*y)); } // G4cout << "e= " << GammaEnergy << " Z= " << Z << " cross= " << xSection << G4endl; return xSection; } //ooo0000000000ooo0000000000 </pre>	G4double y = G4Log(GammaEnergy/10);		G4double y = G4Log(GammaEnergy/I0);
<pre>// G4cout << "e= " << GammaEnergy << " Z= " << Z << " cross= " << xSection << G4endl; return xSection; } // G4cout << "e= " << GammaEnergy << " Z= " << Z // G4cout << "e= " << GammaEnergy << " Z= " << Z // << " cross= " << xSection << G4endl; return xSection; }</pre>	x >ection *= G4Exp(-y*(C1+C2*y));		xsection *= G4Exp(-y*(c1+c2*y));
<pre>return xSection; } if(xSection < 0.0) { xSection = 0.0; } // G4cout << "e= " << GammaEnergy << " Z= " << Z // << " cross= " << xSection << G4endl; return xSection; }</pre>	// G4cout << "e= " << GammaEnergy << " 7= " << 7 << " cross= " << xSection << G4endl:	00]
<pre>} // G4cout << "e= " << GammaEnergy << " Z= " << Z // << " cross= " << xSection << G4endl; return xSection; }</pre>	return xSection;	23	if (xSection < 0.0) { xSection = 0.0; }
//ooo000000000ooo000000000ooo000000	}		// G4cout << "e= " << GammaEnergy << " Z= " << Z
//oooUUUUU0oooooUUUUU0oooooUUUUU0oooooOU0000oo			<pre>// << " cross= " << xSection << G4endl;</pre>
	//ooo000000000		return xSection;

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Beware:

Sometimes automated tools don't catch code duplication effectively

Code reviews!

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Long method

The longer a method is, the more difficult it is to understand

- Decomposing methods
- Most of the time: just Extract Method
 - What to extract?
 - Understand what the code does
 - Comments in the code may help

Large class

- A class that tries to do too much
- Often has too many instance variables

G4VEnergyLossProcess

- Prone to duplicated code
- Extract Class
- Extract SubClass
- Extract Interface

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restance of the second	And the second s		17	<pre>************************************</pre>	A second se	University of the second secon	A second se	Republic to the second	Representation of the second s	A second	A second se	A subscription of the second s
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One class does work that should be done by two

Create a new class and move the relevant fields and methods from the old class into the new class





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Extract interface

Several clients use the same subset of a class's interface, or two classes have part of their interfaces in common

Extract the subset into an abstract interface



Pull up method

Methods with identical results on subclasses





Push down method

Behavior in a superclass is relevant only for some of its subclasses

Move it to those subclasses



Replace inheritance with delegation

A subclass uses only part of a superclass interface or does not want to inherit data

Create a field for the superclass, adjust methods to delegate to the superclass, and remove the subclassing



Refactoring mechanics

Catalog of refactoring techniques

Reference material in Fowler's book and web site

Not meant to be learned by heart, but to be exercised when a pertinent "smell" is identified

Composing Methods

- Extract Method turns a code fragment into a function
- Inline Method is the opposite of Extract Method
- Inline Temp gets rid of a temporary variable by moving the expression to where the temp is used
- Replace Temp with Query removes a temporary variable and instead uses a function call where the temp was used
- Introduce Explaining Variable replaces comments and complex expressions with a temp variable that is well named
- Split Temporary Variable splits a temp that is used for two different things into two different variables
- Remove Assignments to Parameters removes assignments to function parameters within the function
- Replace Method with Method Object moves a complex function to its own class
- Substitute Algorithm

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Moving Features Between Objects

- Move Method moves a method from one class to a more appropriate class
- Move Field moves a field (member object) from one class to another class
- Extract Class pulls a set of methods and fields from one class into a new class
- Inline Class opposite of Extract Class
- Hide Delegate: a class that provides access to an object of another class instead provides the methods of that class by delegation
- **Remove Middle Man** opposite of Hide Delegate
- Introduce Foreign Method adds a method to an untouchable class by passing an instance of the untouchable class into the method
- Introduce Local Extension adds methods to an untouchable class by deriving from it or by wrapping it

Organizing Data

- Self Encapsulate Field creates accessors for private member objects
- Replace Data Value with Object turns a member object into a fullfledged class
- Change Value to Reference
- Change Reference to Value
- Replace Array with Object converts an array which has various fields in each entry into an object
- Duplicate Observed Data introduces a Document/View architecture into an interactive application
- Change Unidirectional Association to Bidirectional introduces a back pointer
- Change Bidirectional Association to Unidirectional is the opposite
- Replace Magic Number with Symbolic Constant

Organizing Data

- Encapsulate Field adds getters and setters
- Encapsulate Collection hides a collection within a class
- Replace Record with Data Class makes a dumb data class to represent a record structure
- Replace Type Code with Class replaces an enumeration type with a class that has a set of global instances of itself, one for each possible value
- Replace Type Code with Subclasses introduces a polymorphic hierarchy to replace an enumeration
- **Replace Type Code with State/Strategy** allows change at runtime
- Replace Subclass with Fields is used when the subclasses no longer serve any real purpose

Simplifying Conditional Expressions

- Decompose Conditional extracts the condition, the "then" part, and the "else" part into functions
- Consolidate Conditional Expression combines a series of "if" into one
- Consolidate Duplicate Conditional Fragments factors out code that is common to a "then" part and an "else" part
- Remove Control Flag replaces flags that trigger exits with return, continue, and break
- Replace Nested Conditional with Guard Clauses replaces nested "if" with returns
- Replace Conditional with Polymorphism replaces case statements with polymorphism
- Introduce Null Object replaces checks for null values with an object
- Introduce Assertion uses assertions to describe a function's preconditions

Making Method Calls Simpler

- Rename Method
- Add Parameter to a function
- Remove Parameter from a function
- Separate Query from Modifier to avoid side-effects
- Parametrize Method reduces a set of similar functions to a single function with a parameter to differentiate amongst the functions
- Replace Parameter with Explicit Methods
- Preserve Whole Object passes an object to a method instead of selected fields
- Replace Parameter with Method reduces a parameter list by using a value that is already available within the class
- Introduce Parameter Object groups parameters into a single object
- **Remove Setting Method** makes an attribute read-only
- Hide Method makes a method private
- Replace Constructor with Factory Method supports polymorphism
 Encapsulate Downcast hides a downcast within a method
- Replace Error Code with Exception separates error-handling from normal paths
- **Replace Exception with Test** provides a method for caller to avoid an exception

Dealing with Generalization

- Pull Up Field factors a common field into a superclass
- Pull Up Method factors a common method into a superclass
- **Push Down Method** moves a unique method down into a subclass
- Push Down Field moves a unique field down into a subclass
- Extract Subclass creates a subclass and moves features into it
- Extract Superclass factors common code into a superclass
- Extract Interface promotes decoupling and partitioning of a class's responsibilities by extracting an interface class
- Collapse Hierarchy combines a subclass and a superclass into one
- Form Template Method factors out common behavior into a superclass
- Replace Inheritance with Delegation Replace Delegation with Inheritance

Big Refactorings

Tease Apart Inheritance

- deals with a messy inheritance hierarchy
- Convert Procedural Design to Objects

Separate Domain from Presentation

- moves domain logic out of the UI classes

Extract Hierarchy

- introduces polymorphism to replace complex conditional code

Computational performance

Refactoring may make the code slower

conventional wisdom

Yes, sometimes...

"First do it, then do it right, then do it fast"

Refactoring **prepares the ground** for computational improvement by providing **clean code**

Refactoring and reengineering physics software

Food for thought...

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ABSTRACT

Geant4 is an object-oriented toolkit for the simulation of the passage of particles through matter. Its development was initially motivated by the requirements of physics experiments at high energy hadron colliders under construction in the last decade of the 20th century. Since its release in 1998, it has been exploited in many different applicative fields, including space science, nuclear physics, medical physics and archaeology. Its valuable support to scientific discovery is demonstrated by more than 16000 citations received in the past 25 years, including notable citations for main discoveries in different fields. This accomplishment shows that well designed software plays a key role in enabling scientific advancement. In this paper we discuss the key principles and the innovative decisions at the basis of Geant4, which made it a game changer in high energy physics and related fields, and outline some considerations regarding future directions.

Smells



Duplicated code Long method Large class Long parameter list Shotgun surgery Feature envy Data clumps Switch statement Parallel inheritance hierarchies Lazy class Speculative generality **Temporary field** Comments **Refused bequest** Primitive obsession Message chains Middle man Inappropriate intimacy

em utils::G4VEmPro flucModel :G4VEmFluctuationMode anglModel :G4VEmAngularDistribution name :G4String {readOnly} lowLimit :G4double highLimit :G4double eMinActive :G4double eMaxActive :G4double polarAngleLimit :G4double ndary Threshold :G4d eLPMflag :G4bool lagDeexcitation :G4boo lagEorceBuildTable 'G4boo incalTable (G4boo) calElmSelectors :G4t nSelectors :std::vecto ElementData :G4ElementData* icleChange :G4VParticleCha tionTable :G4PhysicsTable* eDensity Factor :std::vector<G4double>* (readOnly heDensityIdx :std::vector<G4int>* {readOnly} Wanager :G4LossTableManage CurrentCouple :G4MaterialCutsCouple* (readOnly) CurrentElement :G4Element* {readOnly} xsec :std::vector<G4doubl ~G4VEmModel/ Initialise() :void SampleSecondaries() . InitialiseLocal() :void InitialiseForMaterial() :void InitialiseForElement() :void ComputeDEDXPerVolume() :G-CrossSectionPerVolume() :G4double ComputeCrossSectionPerAtom() G4doub ChargeSquareRatio() :G4double GetChargeSquareRatio() :G4double GetParticleCharge() :G4doub StartTracking() :voi Value() G4double MinPrimaryEnergy() :G4doub MinEnergyCut() :G4double -currentModel SetupForMaterial() :void DefineForRegion() :void GetParticleChangeForLoss() :G4ParticleChangeForLoss* GetParticleChangeForGamma() :G4ParticleChangeForGamma ulidPhysicsTable() :void tartTracking() :void MaxSecondaryEnergy() :G4double InitialiseElementSelectors() :v oid GetElementSelectors() :std::v SetElementSelectors() :void StepDolt() :G4VParticleChar ePhysicsTable() :G4bool ComputeDEDX() :G4double CrossSection() :G4double ComputeMeanFreePath() :G4double computeCrossSectionPerAtom() :G4 SelectIsotopeNumber() (G4int SelectRandomAtom() :G4Elemen SelectRandomAtom() :G4Elemen SelectRandomAtomNumber() :G4ir SetParticleChange() :void SetCrossSectionTable() :void GetElementData() :G4ElementData* GetCrossSectionTable() :G4PhysicsTable* GetModelOfFluctuations() :G4VEmFluctuationMod GetAngularDistribution() :G4VEmAngularDistribution Particle() :G4Particle SetAngularDistribution() :void HighEnergy Limit() :G4double {query} LowEnergy Limit() :G4double {query} HighEnergy Activ ationLimit() :G4doub LowEnergy ActivationLimit() :G4double (query) PolarAngleLimit() :G4double (query) Secondary Threshold() :G4double (query LPMFlag() :G4bool (query) DeexcitationFlag() :G4bool (query) ForceBuildTableFlag() :G4bool (query) UseAngularGeneratorFlag() :G4bool (query) SetAngularGeneratorFlag() :void SetHighEnergyLimit() :void SetLowEnergyLimit() :void SetActivationHighEnergyLimit() :vo SetActivationLowEnergyLimit() :void IsActive() :G4bool SetPolarAngleLimit() :void tor() :G4PI SetSecondary Threshold() :voi SetLPMFlag() :void SetDeexcitationFlag() :void SetForceBuildTable() :void SetMasterThread() :v oid IsMaster() :G4bool (query) MaxSecondary KinEnergy () :G4double GetName() :G4String& (query) SetCurrentCounle() :v oid GetCurrentElement() :G4Element* {query} CurrentCouple() :G4MaterialCutsCouple* {query} SetCurrentElement() :void perator=() :G4VEmModel & GetCurrentLambda() :G4c G4VEmProcess() operator=() :G4VEmProcess 8 G4PEEffectFluoModel theGamma :G4ParticleDefinition theElectron :G4ParticleDefinition* fParticleChange :G4ParticleChangeForGamma fAtom Deexcitation :G4 VAtom Deexcitation* fminimalEnergy :G4double fSandiaCof :std::vector<G4double> G4PhotoElectricEffect G4PEEffectEluoModel() ~G4PEEffectEluoModel(isInitialised :G4bool Initialise() :void ComputeCrossSectionPerAtom() :G4double G4PhotoElectricEffect() ~G4PhotoElectricEffect CrossSectionPerVolume() :G4double SampleSecondaries() :void IsApplicable() :G4bool PrintInfo() :void operator=() :G4PEEffectFluoModel &

InitialiseProcess() :voi

G4PEEffectEluoModel()

em_utils::G4VEmModel

Evolution away from **RD44** discipline

Electromagnetic smells

Coupling

 σ_{tot} and final state modeling have been decoupled in hadronic physics design since RD44

"model"

Total cross section

Whether a process occurs

Final state generation

How a process occurs

Dependencies

on other parts of the software

One needs a geométry (a full-scale simulation applicatio to test a cross section often Difficult to test \rightarrow no testing

Problem domain analysis, improve domain decomposition and software design 71

Maria Grazia Pia, INFN Genova

Magic number

// G4HadronElastic
// Author : [...] 29 June 2009 (redesign old elastic model)

```
G4double dd = 10.;
G4Pow<sup>*</sup> g4pow = G4Pow::GetInstance();
if (A <= 62) {
bb = 14.5 * g4pow -> Z23(A);
aa = g4pow - powZ(A, 1.63)/bb;
cc = 1.4*g4pow->Z13(A)/dd;
} else {
bb = 60.*g4pow->Z13(A);
aa = g4pow - powZ(A, 1.33)/bb;
cc = 0.4*g4pow->powZ(A, 0.4)/dd;
G4double q1 = 1.0 - \text{std::exp}(-bb*tmax);
G4double q2 = 1.0 - std::exp(-dd*tmax);
G4double s1 = q1*aa;
G4double s2 = q2*cc;
```
M. Fowler, *Refactoring*

Number one in the stink parade is **duplicated**

physics

Two Geant4 physics models, identical underlying physics content *(it used to be different)*



Same compatibility with experiment

Penelope

 0.38 ± 0.06

EPDL97

Code bloat

"Livermore"

EPDL97

 0.38 ± 0.06

Burden on

- Software design
- Maintenance
- User support

Unnecessary complexity

Maria Grazia Pia, INFN Genova

M. Fowler, *Refactoring*

Number one in the stink parade is **duplicated code numbers**

Geant 4 Atomic binding energies

IEEE TRANSACTIONS ON NUCLEAR SCIENCE, VOL. 58, NO. 6, DECEMBER 2

Source of epistemic uncertainties?

Evaluation of Atomic Electron Binding Energies for Monte Carlo Particle Transport

Maria Grazia Pia, Hee Seo, Matej Batic, Marcia Begalli, Chan Hyeong Kim, Lina Quintieri, and Paolo Saracco

- 1. Bearden & Burr (1967)
- 2. Carlson
- 3. EADL
- 4. Sevier
- 5. Tol 1978 (Shirley)
- 6. Tol 1996 (Larkins)
- 7. Williams



Maria Grazia Pia, INFN Genova



Benefits Transparency Ease of maintenance Simplicity of testing for V&V

Basic physics V&V can be performed by means of lightweight unit tests

Exploring new physics models is made easier Quantification of accuracy and performance is facilitated

An example of reengineering + new functionality Photon elastic scattering simulation

IEEE TRANSACTIONS ON NUCLEAR SCIENCE, VOL. 59, NO. 4, AUGUST 2012

State of the art

Photon Elastic Scattering Simulation: Validation and Improvements to Geant4

Matej Batič, Gabriela Hoff, Maria Grazia Pia, and Paolo Saracco

Form factor approximation:

Non-relativistic, relativistic, modified + anomalous scattering factors

2nd order S-matrix calculations

recent calculations, not yet used in Monte Carlo codes

Quantification

Statistical analysis, GoF + categorical

	in Gean	t4 l	Differential cross sections					new		
	Penelope Penelope		EPDL	Relativ.	Non-Rel.	Modified	MFF	RFF	SM	
	2001	2008		FF	FF	FF	ASF	ASF	NI	
3	0.27	0.38	0.38	0.25	0.35	0.49	0.52	0.48	0.77	
error	± 0.05	± 0.06	± 0.06	± 0.05	± 0.06	± 0.06	± 0.06	± 0.06	± 0.05	

 ϵ = fraction of test cases not incompatible with experiment, 0.01 significance

Computational performance

Popular belief

Physics model X is intrinsically slow

Baroque methods to combine it with "faster" lower precision models and limit its use to cases where one is willing to pay for higher precision

This design introduces an additional computational burden due to the effects of inheritance and the combination algorithms themselves

Truth

Physics model X is intrinsically fast

But its computationally fast physics functionality is spoiled by an inefficient sampling algorithm



e time

 No code smell
Spotted through in-depth code review in the course of software validation

Maria Grazia Pia, INFN Genova

Change the sampling algorithm! 77

It works It doesn't hurt... so long as you don't want to change it

Most of the fear involved in making changes to large software systems is **fear of introducing bugs**

The legacy code dilemma

When we change code, we should have tests in place

To put tests in place, we often need to change code

With tests, you can change (and improve) your code

Without tests, you just don't know whether things are getting better or worse

Dealing with legacy code

Legacy code often lacks tests

Techniques to make existing code testable

Strategy

- 1. Identify change points
- 2. Find an inflection point
- 3. Cover the inflection point
 - a. Break external dependencies
 - b. Break internal dependencies
 - c. Write tests
- 4. Make changes
- 5. Refactor the covered code



Prevention

- Most of these problems can be easily solved if we simply write tests as we develop our code (and we maintain them...)
- If a test is hard to write, that means that we have to devise a different design which is testable

It is always possible

Provocative thought...

Need to refactor legacy code due to:

- Requirements change
- Computing environment changes (compilers, language standards...)
- New technology becomes available

Need to refactor legacy code due to:

- Laziness to adopt a sound software development process
- Sloppiness
- Refuse to invest in learning technology
- Contempt for good practices
- Lack of design and code reviews
- Lack of adequate mentoring

. . .



Martin Fowler et al. **Refactoring**: Improving the Design of Existing Code Addison-Wesley, 1999/2018 <u>http://www.refactoring.com/</u>

Books





Serge Demeyer, Stéphane Ducasse, Oscar Nierstrasz **Object Oriented Reengineering Patterns** Morgan-Kaufmann, 2003 Revised 2008: http://www.iam.unibe.ch/~scg/OORP



Micheal Feathers **Working Effectively with Legacy Code** Prentice Hall, 2005

Maria Grazia Pia, INFN Genova

Other useful books





Methods

Techniques

Dealing with legacy code in a disciplined, effective way

Sources for further learning

Keep in mind peculiarities of physics software!

Refactoring techniques and reengineering patterns support software maintenance, evolution and validation They may also contribute to improving **computational performance**

Thorough testing is the key

...but also sound background in OO methods, healthy software engineering practices and physics insight

Conclusions



Hands-on exercises

"Video store" example

- M. Fowler, Refactoring, chapter 1
- Simple problem domain
- Practice a few refactoring techniques
- Learn the method
 - Write tests, change in small steps, test, change, test...
- We'll figure out together how many steps we do as an exercise and how many we'll just review