

**Attribution of Responsibility  
for Climate Change**

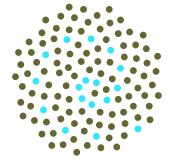
**The Mathematics  
Behind the Brazilian Proposal**

Ian G. Enting

MASCOS

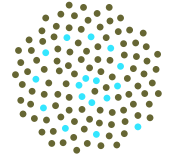
The University of Melbourne

# Acknowledgments



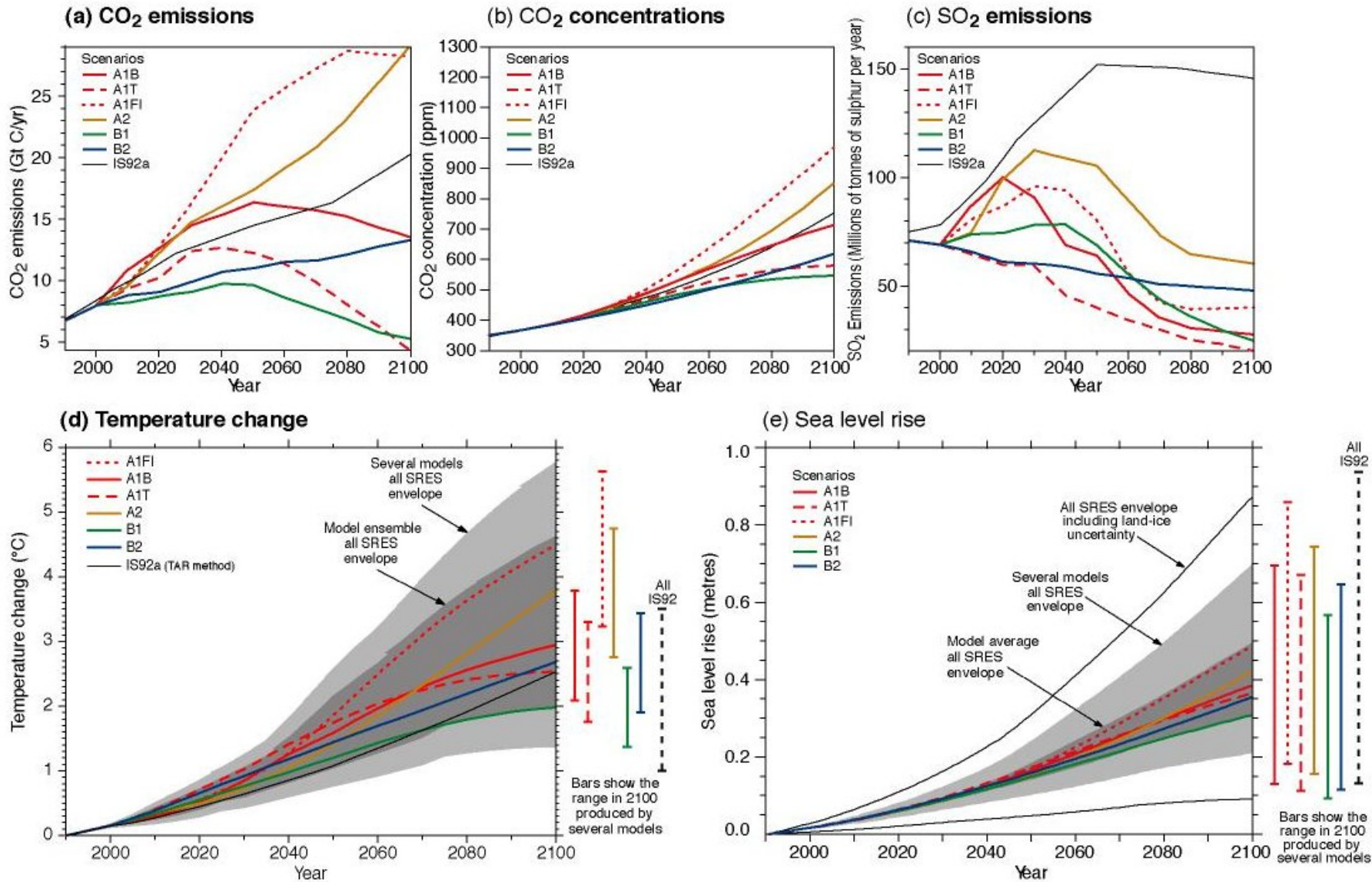
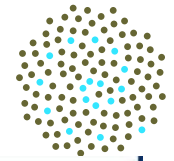
- The Center of Excellence for Mathematics and Statistics of Complex Systems (MASCOS) is funded by the Australian Research Council (ARC).
- My fellowship at MASCOS is supported by CSIRO through a sponsorship agreement.
- Collaborators: Cathy Trudinger of CSIRO Marine and Atmospheric Research and members of the MATCH working group on the Brazilian Proposal.

# Summary



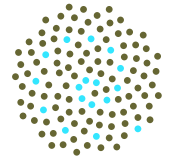
- Time-scales for the greenhouse effect
  - committed warming
- The Brazilian Proposal
  - setting reduction targets in proportion to responsibility
- Adjoint modelling & automatic differentiation
  - efficient calculation of sensitivities
  - operator overloading
- Analysing the Brazilian proposal
  - Who's to blame for the greenhouse effect?

# The causal chain



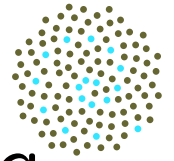
From IPCC Third Assessment report

# IPCC is Cautious

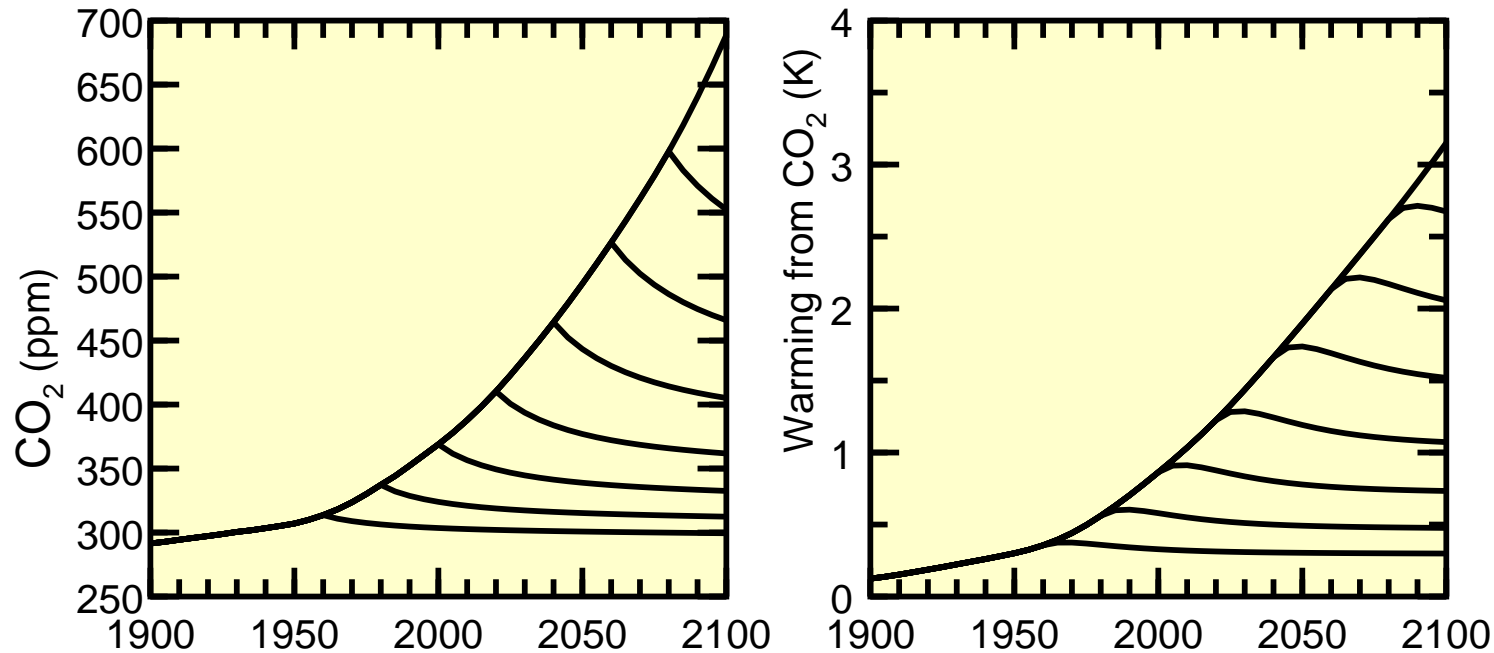


- The IPCC requirement for a consensus evaluation of ‘well-established science’ can mean that IPCC reports lag behind the forefront of science, sometimes by more than a decade.
- Similarly, Al Gore’s book and film are very careful to avoid premature speculation about extreme possibilities.
- Recent study (Science Express, 1/2/07) indicates that changes since 2001 are tracking the high end of IPCC projections.
- Overview by Barrie Pittock on mechanisms that may imply current under-estimation of change.
- For ‘alarmist’ views, see ‘Pentagon scenarios’.

# Timescales



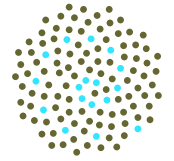
CO<sub>2</sub> concentrations and consequent warming, partitioned according to time of emission.



Lowest bands are from pre-1960 emissions, next from 1960 to 1980 emissions, etc.

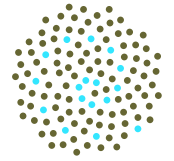
Increase in contribution to warming after time of emissions from 'committed warming' effect.

# Some key dates



- circa 1960: Keeling measures CO<sub>2</sub> growth
- Oct 1985. *it is now believed that in the first half of the next century a rise of global mean temperature could occur which is greater than any in man's history.*  
(UNEP/WMO/ICSU conference, Villach).
- 1998. IPCC established by WMO and UNEP
- UNFCCC. Rio 1992. In force 21/3/99
- 1995. CoP-1, Berlin Mandate (to negotiate a protocol)
- Kyoto Protocol: 11/12/97. came into force 16/2/05.
- 2001 (IPCC) balance of evidence suggests discernable human influence
- Feb 2007 (IPCC): **unequivocal** attribution to human agency of most of the warming since 1950.
- 2008-2012. Kyoto Protocol: first commitment period

# Mitigation frameworks

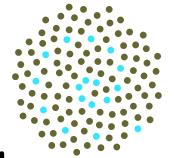


- Kyoto, specified percentage reductions
- Contraction and convergence
- Sector-based targets – e.g. the Tryptique framework underlying target-setting within the ‘EU bubble’

Note that measures such as carbon trading, clean development mechanism or the McKibbin-Wilcoxon multi-level permits are not primarily about mitigation levels, they are about economically-efficient implementation.



# Brazilian Proposal



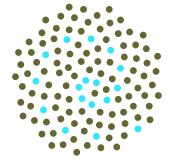
Tabled by Brazil during negotiations leading to Kyoto Protocol — Flicked-passed to Subsidiary Body for Scientific and Technical Advice (SBSTA).

Proposes that emission reduction targets should be proportional to nations' relative responsibility for the greenhouse effect.

Issues:

- Indicator? What quantity is used as a measure of the greenhouse effect?
- For what period of emissions is responsibility attributed?
- How are non-linear responses attributed?

# Political significance

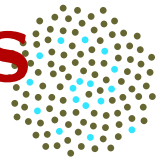


- It captures the historical responsibility of developed nations (for global warming) and the requirement of developed nations to take the lead in combating climate change — i.e. as prescribed by the UN Framework Convention on Climate Change.
- It provides a formula that can apply to all nations
- Therefore it provides a way of engaging developing nations as they develop.

# Expert Working Group: MATCH

- Initially under the auspices of SBSTA, initial meeting in Brazil
- Now an informal working group with government support from UK, FRG, Norway. Expert Meetings: Bonn, 2001; Bracknell, UK, 2002; Berlin, 2003; Cologne, 2004; Rio de Janeiro, 2005; Reading, UK 2005; Louvain-la-Neuve, Belgium 2006; Cologne, 2006.
- Joint paper published:  
den Elzen et al. Analysing countries contributions to climate change: scientific and policy-related choices. *Environmental Science and Policy* 8: 614–635 (2005).
- Initial results presented to SBSTA, May 2006. Final results scheduled for presentation to SBSTA late 2007.

# Brazilian Proposal as Derivatives



As example, use indicator  $T^* = T_{\text{CO}_2}(2000) =$  warming in 2000 from  $\text{CO}_2$  emissions.

$T^*$  is to be attributed to emissions  $E_j(t)$  from country  $j$  with  $E(t) = \sum_j E_j(t)$ .

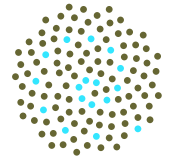
Differential attribution to country  $j$  of emissions at time  $t$  is

$$\frac{\partial T^*}{\partial E_j(t)} E_j(t) = \frac{\partial T^*}{\partial E(t)} E_j(t) = S(t) E_j(t)$$

where  $S(t)$  is a Fréchet derivative.

Cumulated attribution:  $T_j^* = \int S(t) E_j(t) dt$

# Aims of adjoint modelling



Aim is to simplify calculations by separating parametric differentiation from model integration, expressed here in terms of Green's function  $\mathcal{G}$  of  $\mathcal{L}\underline{u}(\cdot) = \underline{f}(\cdot)[\underline{a}]$ .

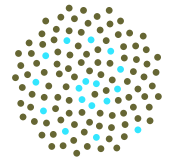
Considers  $\nabla_{\underline{a}}\langle \underline{w}(\cdot) | \underline{u}(\cdot)[\underline{a}] \rangle$   
where  $\underline{u}(\cdot)[\underline{a}] = \mathcal{G}\underline{f}(\cdot)[\underline{a}]$  (or  $\mathcal{L}\underline{u}(\cdot) = \underline{f}(\cdot)[\underline{a}]$ )

Transforms as

$$\nabla_{\underline{a}}\langle \underline{w}(\cdot) | \underline{u}(\cdot)[\underline{a}] \rangle = \nabla_{\underline{a}}\langle \underline{w}(\cdot) | \mathcal{G}\underline{f}(\cdot)[\underline{a}] \rangle =$$
$$\nabla_{\underline{a}}\langle \mathcal{G}^\dagger \underline{w}(\cdot) | \underline{f}(\cdot)[\underline{a}] \rangle = \nabla_{\underline{a}}\langle \underline{v}(\cdot) | \underline{f}(\cdot)[\underline{a}] \rangle$$

where  $\underline{v}(\cdot) = \mathcal{G}^\dagger \underline{w}(\cdot)$  defines a single function  $\underline{v}(\cdot)$  with no dependence on  $\underline{a}$ .

# Principles of adjoint modelling



Given  $\underline{u}(\cdot)[\underline{a}] = \mathcal{G}\underline{f}(\cdot)[\underline{a}]$ , where typically  $\mathcal{L}\underline{u}(\cdot)$  is linearisation of a more general model:

Formally:

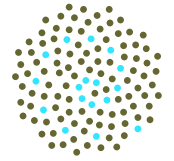
$$\begin{aligned}\nabla_{\underline{a}}\langle\underline{w}(\cdot)|\underline{u}(\cdot)[\underline{a}]\rangle &= \nabla_{\underline{a}}\langle\underline{w}(\cdot)|\mathcal{G}\underline{f}(\cdot)[\underline{a}]\rangle = \\ \nabla_{\underline{a}}\langle\mathcal{G}^\dagger\underline{w}(\cdot)|\underline{f}(\cdot)[\underline{a}]\rangle &= \nabla_{\underline{a}}\langle\underline{v}(\cdot)|\underline{f}(\cdot)[\underline{a}]\rangle \\ \text{with } \underline{v}(\cdot) &= \mathcal{G}^\dagger\underline{w}(\cdot)\end{aligned}$$

In practice, used as  $\mathcal{L}\underline{u}(\cdot) = \underline{f}(\cdot)[\underline{a}]$

$$\begin{aligned}\nabla_{\underline{a}}\langle\underline{w}(\cdot)|\underline{u}(\cdot)[\underline{a}]\rangle &= \nabla_{\underline{a}}\langle\mathcal{L}^\dagger\underline{v}(\cdot)|\underline{u}(\cdot)[\underline{a}]\rangle = \\ \nabla_{\underline{a}}\langle\underline{v}(\cdot)|\mathcal{L}\underline{u}(\cdot)[\underline{a}]\rangle &= \nabla_{\underline{a}}\langle\underline{v}(\cdot)|\underline{f}(\cdot)[\underline{a}]\rangle\end{aligned}$$

with  $\underline{w}(\cdot) = \mathcal{L}^\dagger\underline{v}(\cdot)$  giving equations for adjoint model.

# Applying adjoint modelling



**Differentiation** (only case used in this talk)

$$\nabla_{\underline{a}} \langle \underline{w}(\cdot) | \mathcal{G} \underline{f}(\cdot) [\underline{a}] \rangle = \nabla_{\underline{a}} \langle \mathcal{G}^\dagger \underline{w}(\cdot) | \underline{f}(\cdot) [\underline{a}] \rangle$$

**Gradients for soft constraints .**

$$\begin{aligned} \nabla_{\underline{a}} \langle \underline{H} \underline{u} - \underline{z} | \underline{H} \underline{u} - \underline{z} \rangle &= 2 \nabla_{\underline{a}} \langle \underline{H} \underline{u}_0 - \underline{z} | \underline{H} \underline{u} \rangle = \\ 2 \nabla_{\underline{a}} \langle \underline{H} \underline{u}_0 - \underline{z} | \underline{H} \mathcal{L} \underline{f} \rangle &= 2 \nabla_{\underline{a}} \langle (\underline{H} \mathcal{L})^\dagger (\underline{H} \underline{u}_0 - \underline{z}) | \underline{f} \rangle \end{aligned}$$

**Gradients, with hard constraints:  $\mathcal{L} \underline{u}(\cdot) = 0$**

$$\Theta^* = \Theta - \langle \underline{v}(\cdot) | \mathcal{L} \underline{u}(\cdot) \rangle$$

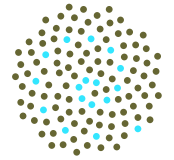
The function  $\underline{v}(\cdot)$  is the Lagrange multiplier.

$$\nabla_{\underline{u}} \Theta^* = \nabla_{\underline{u}} \Theta - \nabla_{\underline{u}} \langle \mathcal{L}^\dagger | \underline{u}(\cdot) \rangle, \text{ whence}$$

$$\mathcal{L}^\dagger \underline{v}(\cdot) = \nabla_{\underline{u}} \Theta \text{ — the adjoint equations}$$

define the Lagrange multiplier

# Tangent Linear Model (TLM)



For  $N$  DEs:  $\frac{d}{dt}x_j = g_j(\{x_k\}, a, t)$  for  $j = 1, N$

we can define sensitivities as

$$y_j = \frac{\partial}{\partial a} x_j \quad \text{for } j = 1, N \quad \text{or} \quad y_{j,p} = \frac{\partial}{\partial a_p} x_j$$

to give 'tangent linear model(s)':

$$\frac{d}{dt}y_m = \frac{\partial}{\partial a} g_m(\{x_k\}, a, t) + \sum_n \frac{\partial}{\partial x_n} g_m(\{x_k\}, a, t) y_n$$

$$\frac{d}{dt}y_{m,p} = \frac{\partial}{\partial a_p} g_m(\{x_k\}, \underline{a}, t) + \sum_n \frac{\partial}{\partial x_n} g_m(\{x_k\}, \underline{a}, t) y_{n,p}$$



# Algorithmic Differentiation (AD)

Differentiation by successive use of chain rule.

For binary operation  $c = f(a, b)$ ,

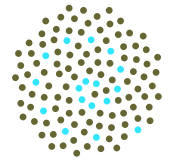
$$\frac{\partial c}{\partial \alpha} = \frac{\partial f}{\partial a} * \frac{\partial a}{\partial \alpha} + \frac{\partial f}{\partial b} * \frac{\partial b}{\partial \alpha}$$

e.g.  $c = a + b \quad \rightarrow \quad \frac{\partial c}{\partial \alpha} = \frac{\partial a}{\partial \alpha} + \frac{\partial b}{\partial \alpha}$

or  $c = a * b \quad \rightarrow \quad \frac{\partial c}{\partial \alpha} = b * \frac{\partial a}{\partial \alpha} + a * \frac{\partial b}{\partial \alpha}$

Automatically derives TLM: converting program to code for derivatives, one operation at a time.

# Approaches to AD



Hand-code program to calculate derivatives — laborious, error-prone and must be repeated each time the model changes.

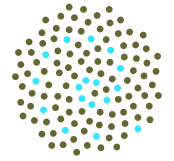
Symbolic algebra (e.g. Mathematica) — problematic for adjoints.

Tangent/adjoint compilers — transform source into code for tangent or adjoint models.

Operator overloading to produce a ‘script’ that is analysed to give code for the derivatives.

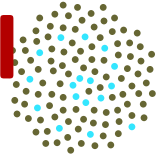
Use operator overloading capabilities directly — straightforward for tangent-linear-model, but restricted applicability to adjoint models.

# AD by operator overloading



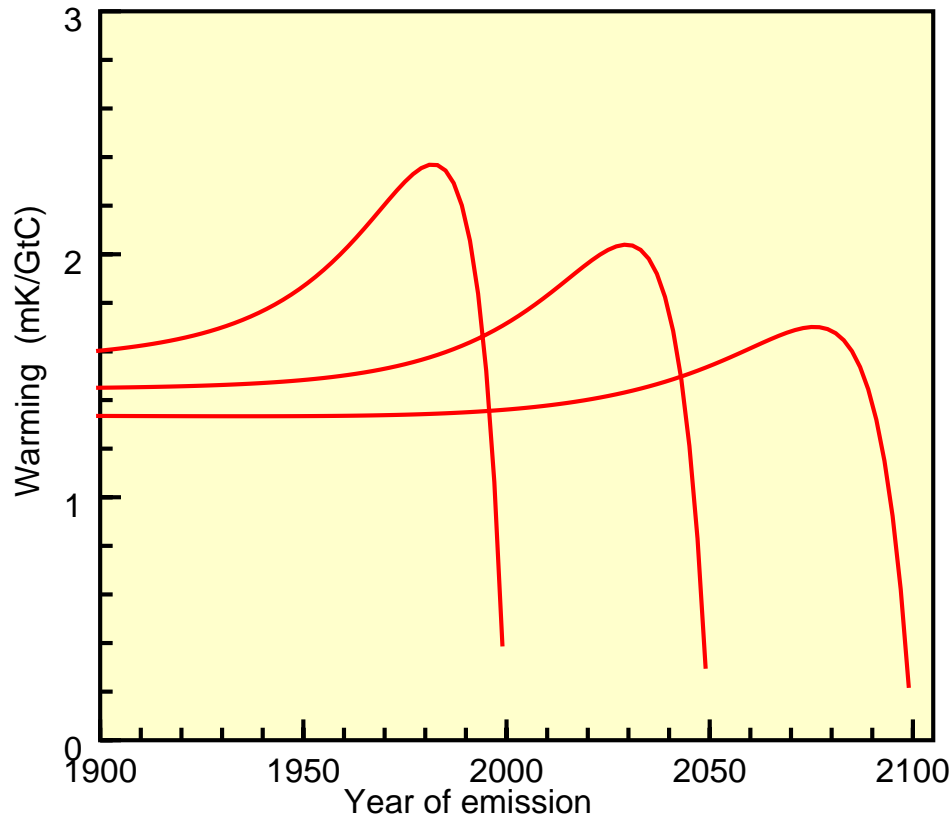
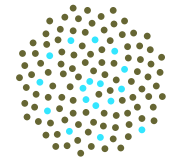
- Needs object-oriented language, C++ or F90
- Define new types that include derivative information
- Use overloading to define operations on these types
- Modify type declarations in original model code to invoke new types
- Set up requisite I/O for derivative information

# Analysing the Brazilian Proposal



- Construct simple climate model
- Construct linearisation (e.g. by automatic differentiation)
- Calculate sensitivities, either by brute force application of linearised model or by explicit adjoint model.
- Apply sensitivities to histories of emissions from each nation
- Repeat for all greenhouse gases

# Results: Fréchet Derivatives

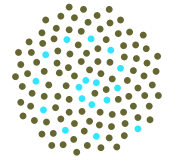


Assumes IS92a emissions. Represents temperature by response function. Linear responses for ocean and biotic carbon, coupled non-linearly to atmospheric CO<sub>2</sub> (as in CSIRO study).

$$\frac{\partial}{\partial E(t)} T(\tau) \text{ for } \tau = 2000, 2050, 2100.$$

Decrease as  $t \rightarrow \tau$  shows 'committed warming'. At any time, warming from the most recent releases is yet to happen.

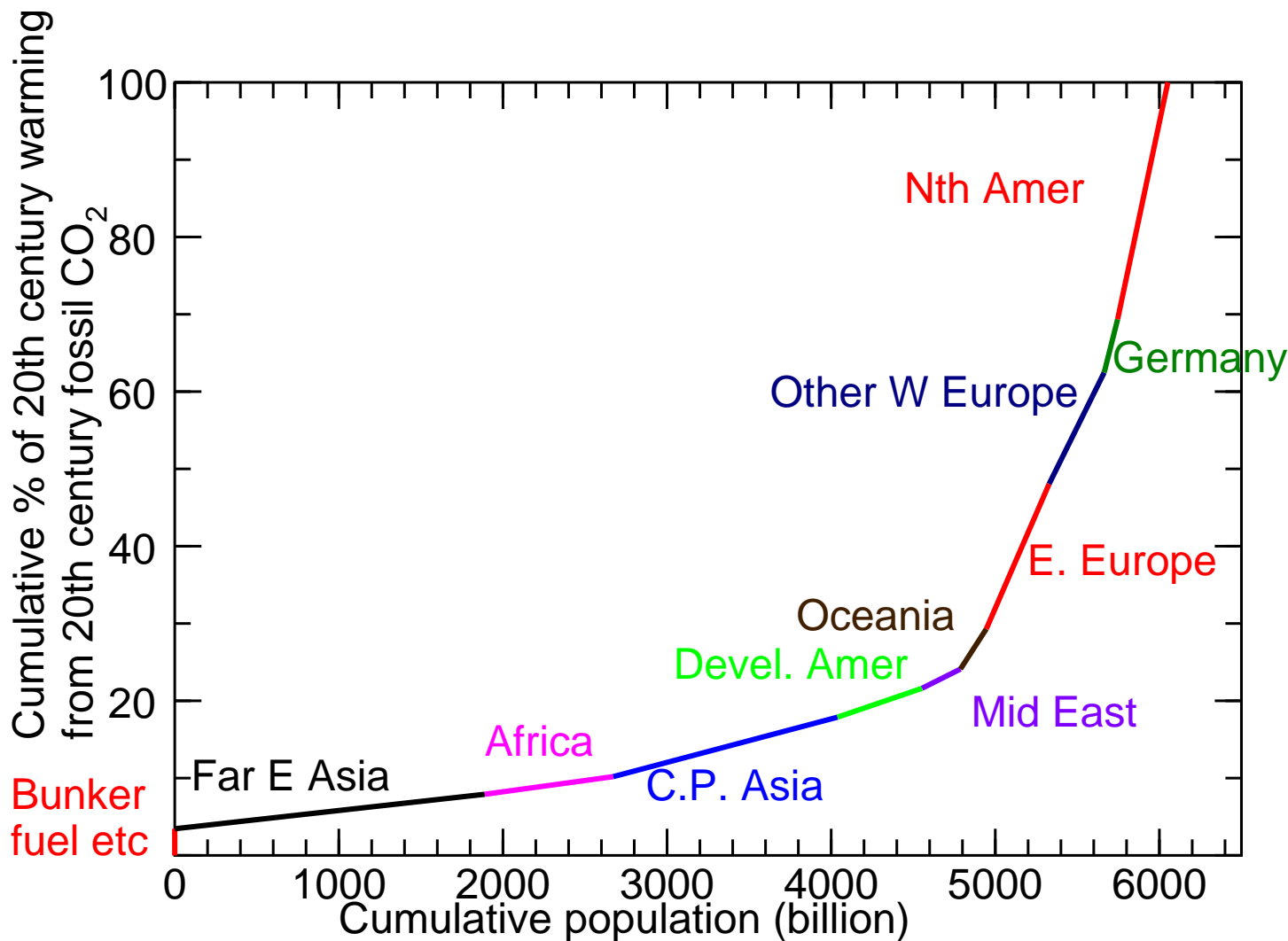
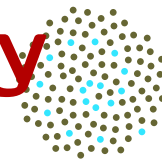
# Implications



- For a given indicator,  $T^*$ , calculation of  $S(t)$  allows attribution to any nation.
- $S(t)$  most efficiently calculated from adjoint model, but for multiple indicator times, tangent linear model not too inefficient.
- Sensitivity of  $T_j^*$  to model uncertainties can be obtained as second derivatives.
- Sensitivity of  $T_j^*$  to uncertainties in emissions can be obtained as

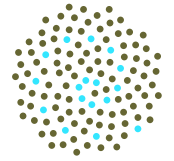
$$\text{Var}[T_j^*] = \int \int S(t) \text{Cov}[E_j(t), E_j(t')] S(t') dt' dt$$

# Attribution 2000, fossil CO<sub>2</sub> only



Cumulative responsibility for the fossil CO<sub>2</sub> component of warming vs cumulative population.

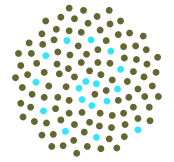
# Concluding remarks



- An interesting example of adjoint sensitivity analysis and automatic differentiation
- The Brazilian Proposal is on the agenda, for formal consideration by Conference of Parties (to the Climate Change Convention) in 2008
- Expert working group is extending calculations to include all major greenhouse gases, with detailed national attribution



# Further Information



MATCH website (Brazilian Proposal):

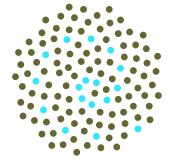
*<http://www/match-info.net>*

Andreas Griewank, 2000, *Evaluating Derivatives: Principles and Techniques of Algorithmic Differentiation*, (SIAM, Philadelphia).

I.G. Enting, 2005, *Automatic differentiation in the analysis of strategies for mitigation of global change*, International Congress on Modelling and Simulation, Melbourne, 2005. Ed. A. Zenger and R. M. Argent, 7pp  
<http://www.mssanz.org.au/modsim05/papers/enting.pdf>

My AD talk from Berkeley is on my website.

# Class Definitions

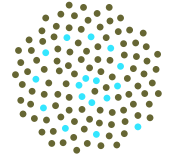


Fragment of C++ class definition to implement operator overloading:

```
class Xvar{
public :
static const int ns = _NUMDERIVS+1;
double xs[_NUMDERIVS+1];
Xvar operator*(Xvar);
...
};

Xvar Xvar::operator*(Xvar b){ Xvar c;
for (int i=1; i < ns; i++)
    c.xs[i] = xs[i]*b.xs[0]+xs[0]*b.xs[i];
c.xs[0] = xs[0]*b.xs[0];
return c;} ;
...
```

# Usage



## Original

```
double F_co2(double c){  
double a;  
a = log(c/280.0)*5.35;  
return a;  
};  
...  
double cc;  
...  
ff = F_CO2(cc)
```

## Transformed

```
Xvar F_co2(Xvar c){  
Xvar a;  
a = log(c/280.0)*5.35;  
return a;  
};  
...  
Xvar cc;  
// Derivatives wrt  
// initial value of cc  
cc.set(280,1);  
...  
ff = F_CO2(cc)
```