



Spectral methods

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MTMW14

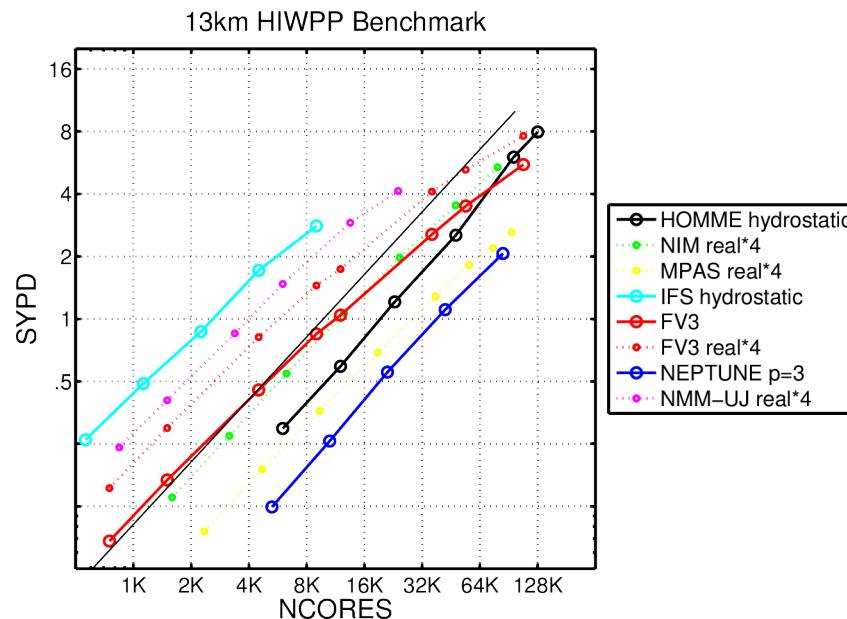
9 March 2023

Based largely on Dave Randall's AT604 lecture notes

Spectral models: very much alive, despite their death having been announced for 20 years.

NGGPS Scaling Study

Spectral models are nowadays fast, efficient, **scalable**, and able to perform scale separation by design.



Thomas Schultess (ETH and CSCS) showed how the IFS can take up the entire Piz Daint (one of the top 10 supercomputers) and scale well at 1km

Nils Wedi carried on to Summit...

<http://www.nws.noaa.gov/ost/nggps/dycoretesting.html>

Recent breakthroughs with the ECMWF IFS

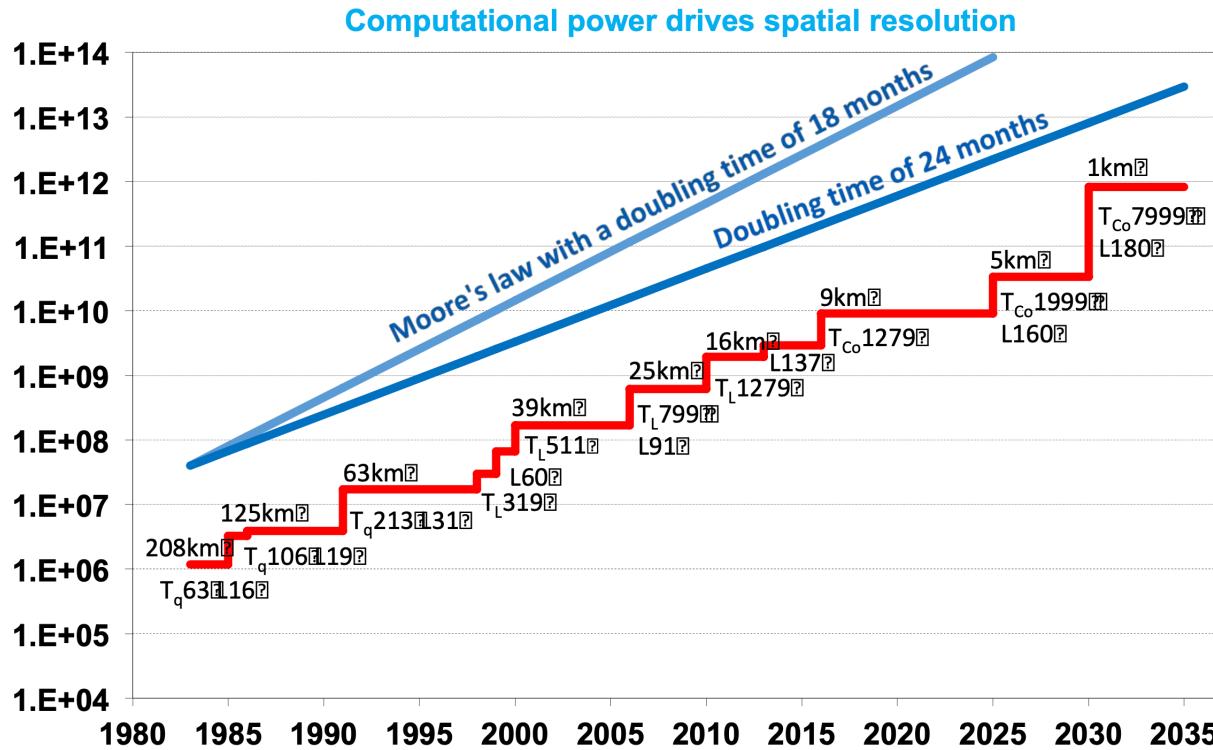


Figure 2: "Understanding grows only logarithmically with the number of floating point operations." (John P. Boyd). The progress in the degrees of freedom (vertical levels x grid columns x prognostic variables) of the ECMWF operational global atmosphere model in comparison to Moore's law, adjusted by ($P^{3/4}$) to account for the change in time-resolution also required in a 3-dimensional dynamical system when increasing spatial resolution. ECMWF's actual progress up to 2018 doubles performance every 24 month. As illustrated, the ambitious goal of reaching a 1 km horizontal resolution with 180 levels in 2030 requires a faster progress. The numbers indicate the average grid-point distance in kilometers and the corresponding spectral resolution and levels used.

Reflecting on the Goal and Baseline for Exascale Computing: A Roadmap Based on Weather and Climate Simulations

Thomas C. Schultheiss
ETH Zurich, Swiss National Supercomputing Centre

Peter Bauer
European Centre for Medium-Range Weather Forecasts

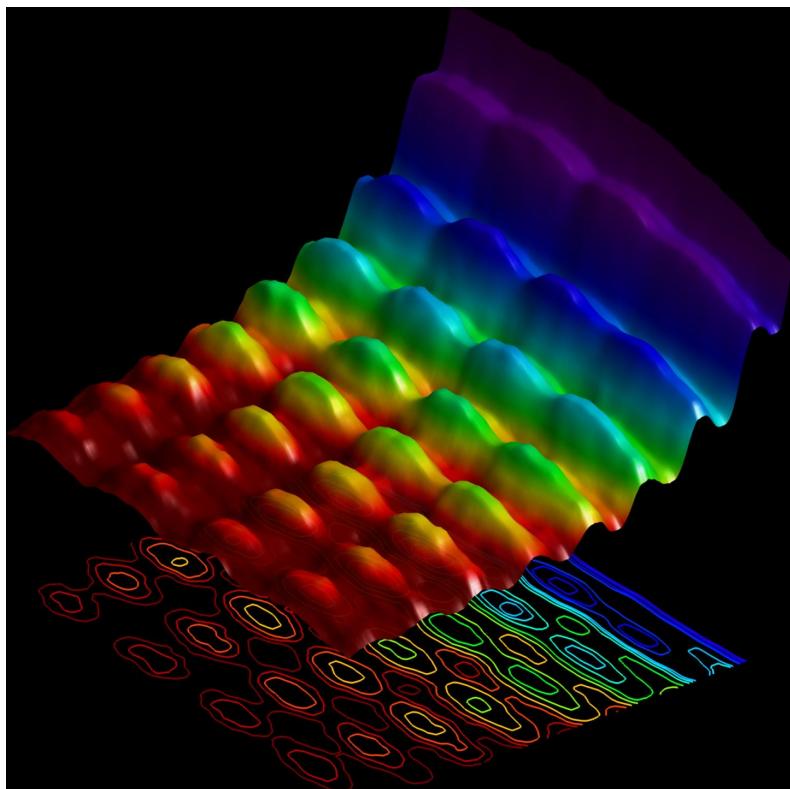
Nils Wedi
European Centre for Medium-Range Weather Forecasts

Oliver Fuhrer
MeteoSwiss
Torsten Hoefer
ETH Zurich
Christoph Schär
ETH Zurich

Table 1. Ambitious target configuration for global weather and climate simulations with km-scale horizontal resolution accounting for physical Earth-system processes, and with today's computational throughput rate.

Horizontal resolution	1 km (globally quasi-uniform)
Vertical resolution	180 levels (surface to ~ 100 km)
Time resolution	0.5 min
Coupled	Land-surface/ocean/ocean-waves/sea-ice
Atmosphere	Non-hydrostatic
Precision	Single or mixed precision
Compute rate	1 SYPD (simulated years per wall-clock day)

Waves or particles?



This photo was published a few days ago in Science. Fabrizio Carbone (EPFL) was able, for the first time, to capture the dual nature of light.

So far we have been dealing in this course with discrete representations of functions, for instance as points or volumes.

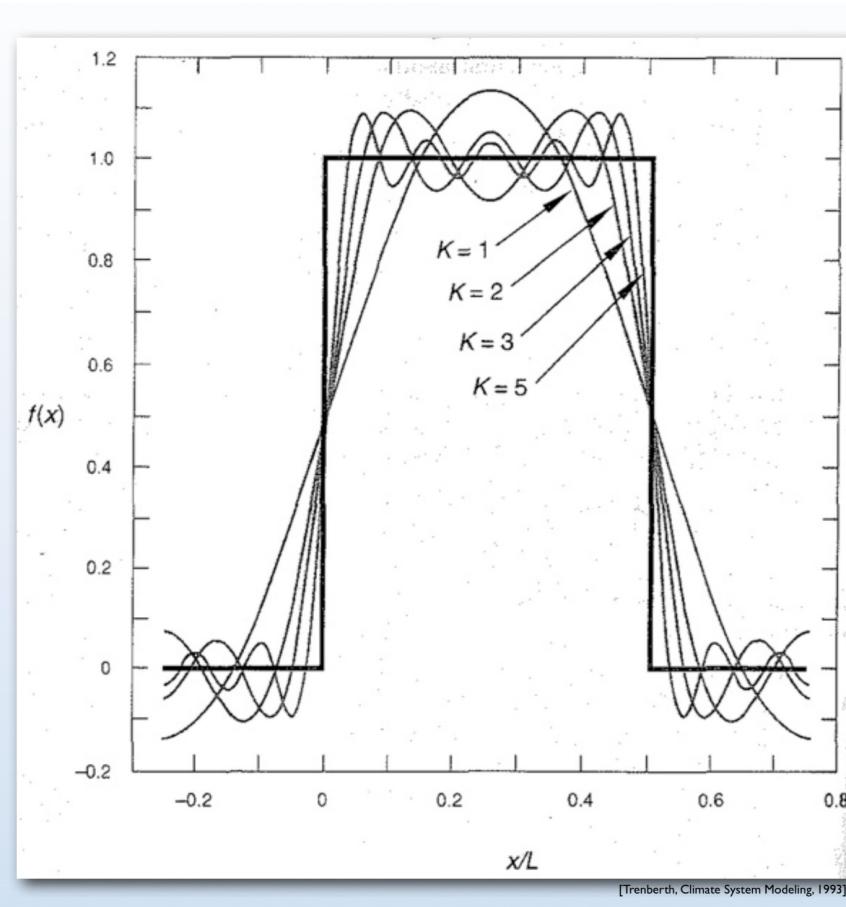
Every continuous function in the function space can be represented as a [linear combination of basis functions](#).

Fourier series are an example of basis functions. From Fourier's theorem, any continuous function can be represented as a sum of *sin* and *cos* functions:

$$q(x, t) = \sum_{k=-\infty}^{\infty} \widehat{q}_k(t) e^{ikx}$$

Back to our square wave example

Improved
Spectral Representation
of Steep Gradients by
Increasing Wave
Numbers



Credit AWI, Germany

In practice, when using a Fourier basis:

- Start with a real and integrable function $q(x,t)$:
- Use the definition:

$$q(x, t) = \sum_{k=-\infty}^{\infty} \widehat{q}_k(t) e^{ikx} \quad (1)$$

Representation of q at point location depends on combination of an infinite number of waves

- Where the complex coefficients can be evaluated:

$$\widehat{q}_k(t) = \frac{1}{L} \int_{x-\frac{L}{2}}^{x+\frac{L}{2}} q(x, t) e^{-ikx} dx \quad (2)$$

Wave characteristics depend on sampling q over a domain of size L

- (1) and (2) form a *transform pair*

Why bother?

- For starters, gradients are exact and easy. For instance, in physical space (x domain):

$$q(x, t) = \sum_{k=-\infty}^{\infty} \hat{q}_k(t) e^{ikx} \quad (1) \quad \rightarrow \quad \frac{\partial q(x, t)}{\partial x} = \sum_{k=-\infty}^{\infty} ik \hat{q}_k(t) e^{ikx}$$

- In practice, for a finite spectral model, the *transform pair* is:

n wave numbers at each grid point j

$$q(x_j, t) \cong \sum_{k=-n}^n \hat{q}_k(t) e^{ikx_j} \quad (6)$$

$$-n \leq k \leq n$$

N is the TRUNCATION

M grid points used to find spectral coefficients

$$\hat{q}_k(t) \cong \frac{1}{M} \sum_{j=1}^M q(x_j, t) e^{-ikx_j} \quad (7)$$

M is the RESOLUTION

$$\frac{\partial q(x_j, t)}{\partial x} \cong \sum_{k=-n}^n ik \hat{q}_k(t) e^{ikx_j}$$

Gradient computed using waves

Truncation, accuracy and computational costs

- The truncation is exactly reversible if M is large. This is equivalent to model resolution in finite difference models
- What are the costs involved? Substitute (6) into (7)

$$\hat{q}_k(t) = \frac{1}{M} \sum_{j=1}^M \left\{ \left[\sum_{l=-n}^n \hat{q}_l(t) e^{ilx_j} \right] e^{-ikx_j} \right\}, \quad -n \leq k \leq n$$
$$-n \leq l \leq n$$

- Both wave numbers bound by n . Max wave number: $\pm 2n$
→ We need $2n+1$ complex conjugates
- The Fourier representation is equivalent to discretising $q(x,t)$ onto **2n+1 grid points**

Waves and particles again

We have:

- Truncation $n \rightarrow$ max wave number: $2n$
- $2n+1$ complex conjugates $\hat{q}_k(t)$

This means that for a grid of M points to represent amplitudes and phases of all waves up to $k = \pm n$, it is necessary that $M \geq 2n + 1$

The truncation in the representation of the waves and the grid resolution are related. This is important, because,

1. as we will see, physical parameterisations operate on the grid, not in spectral space. The model algorithm uses the transform pair to jump from the grid into spectral space and back again.
2. There are non-linear terms...

Equivalence of spectral and finite difference methods

$$\frac{\partial q}{\partial x}(x_l, t) \cong \sum_{k=-n}^n \left[\frac{ik}{M} \sum_{j=1}^M q(x_j, t) e^{-kx_j} \right] e^{ikx_l}$$

Now reverse the two \sum

$$\frac{\partial q}{\partial x}(x_l, t) \cong \sum_{j=1}^M \alpha_j^l q(x_j, t)$$

where $\alpha_j^l \equiv \frac{i}{M} \sum_{k=-n}^n k e^{ik(x_l - x_j)}$

This can be **interpreted as a finite difference approximation.**

A member of the families we have seen before, but now it involves all points in the domain

Let's try to build a simple model: 1D linear advection in spectral space

$$\frac{\partial q}{\partial t} + c \frac{\partial q}{\partial x} = 0$$

$$\sum_{k=-n}^n \frac{d\hat{q}_k}{dt} e^{ikx_j} + c \sum_{k=-n}^n ik\hat{q}_k e^{ikx_j} = 0$$

Since we can invoke linear independence:

$$\frac{d\hat{q}_k}{dt} + ikc\hat{q}_k = 0 \quad \text{for all } -n \leq k \leq n \quad \text{We have our prognostic model!!}$$

remember that $\frac{dq_0}{dt} \equiv 0$

1D linear advection: spectral and finite difference

$$\frac{d\hat{q}_k}{dt} + ikc\hat{q}_k = 0$$

SPECTRAL

$$\frac{d\hat{q}_k}{dt} + ikc \left[\frac{\sin(k\Delta x)}{k\Delta x} \right] \hat{q}_k = 0$$

FINITE DIFFERENCE

Which model will give better results?

More advantages for spectral models: a boundary value problem

$$\nabla^2 q = f(x, y) \quad \text{in 1D: } \frac{d^2 q}{dx^2} = f(x)$$

with periodic B.C.s:

$$\sum_{k=-n}^n (-k^2) \hat{q}_k e^{ikx} = \sum_{k=-n}^n \hat{f}_k e^{ikx}$$

$$\Rightarrow \hat{q}_k = -\frac{f}{k^2} \quad \text{for } -n \leq k \leq n, \text{ not valid for } k = 0$$

This is an exact solution for all modes that are included.

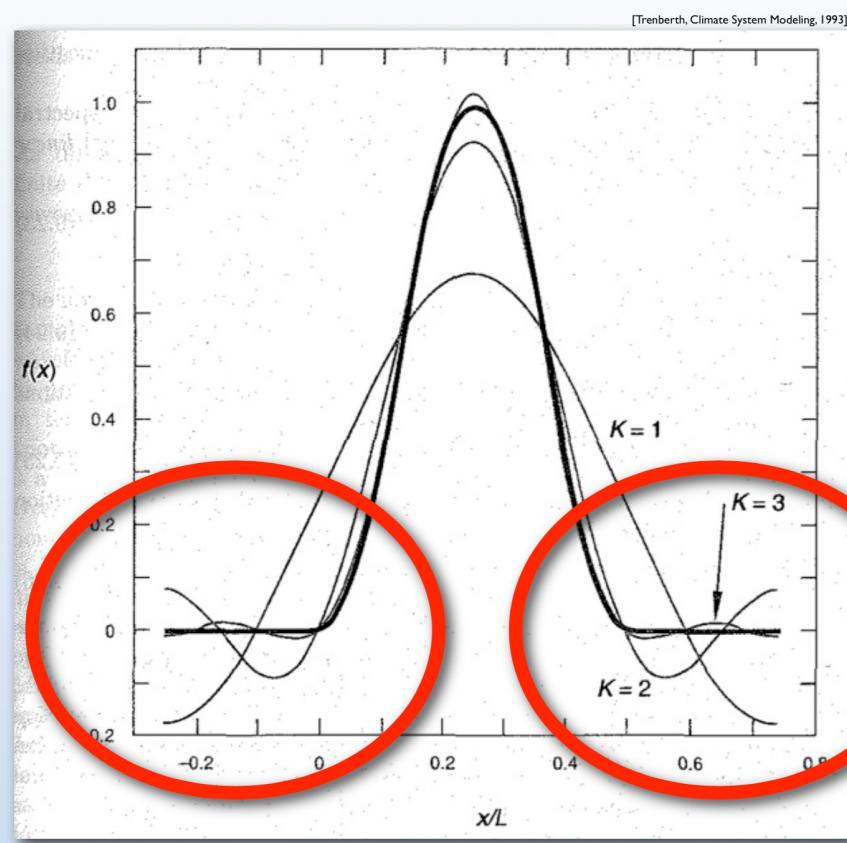
Ends up not being exact because some modes are left out.

Disadvantages I: treatment of orography

This is more typical of older spectral models

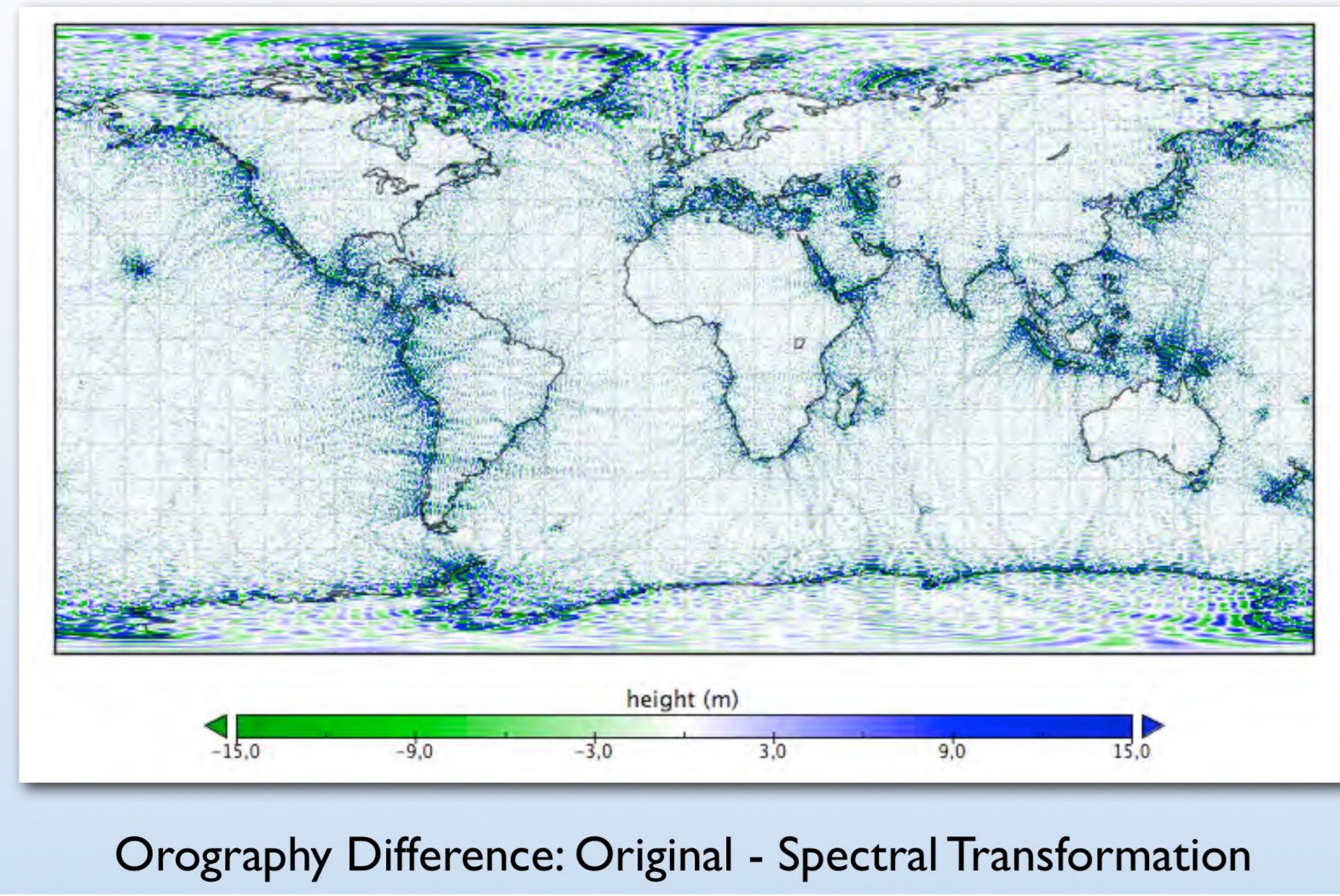
“Overshooting” of a
Spectral Representation
of Steep Gradients

We said that we
can represent
anything, e.g. a
square wave, but
“overshooting” is
one of the many
prices we pay
when we use
spectral methods.



Credit AWI, Germany

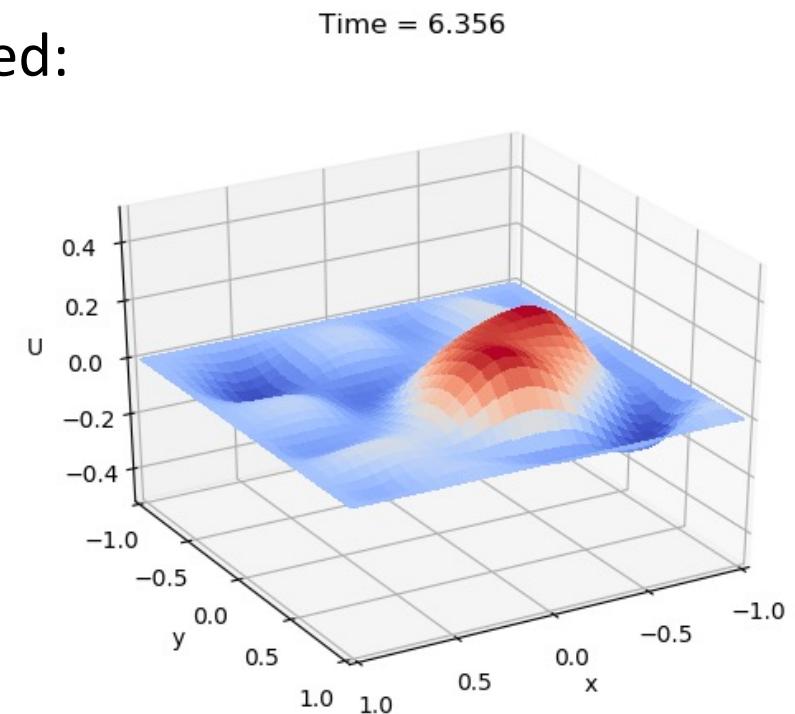
Example: “Overshooting” of Topography Gradients after Transformation into Spectral Space



Brief interval: SWE re-visited with finite differences vs spectral method

For a reasonably smooth initial condition, I used:

	Finite Difference	Spectral	Spectral
Resolution	M=50 grid points in x,y	T=12 wavenumbers	T=24
Time step	0.2s	0.2s	0.2s
Time to solution	30s	15s	59s
But the spectral method converges quickly with low wavenumber, so T=24 is not needed for this problem!!! Also, I am not using any of the typical spectral tricks...			



From python Spectral_wave2_32.py

A bit more trouble when waves interact...

- Let us go back to advection, but this time let us consider the full (non-linear) problem of momentum advection over a periodic domain:

$$\frac{\partial u}{\partial t} + u \frac{\partial u}{\partial x} = 0 \quad \xrightarrow{\hspace{1cm}} \quad \sum_{k=-n}^n \frac{d\hat{u}_k}{dt} e^{ikx} = - \left(\sum_{l=-n}^n \hat{u}_l e^{ilx} \right) \left(\sum_{m=-n}^n i m \hat{u}_m e^{imx} \right)$$

- But now we have products in this form: $e^{ilx} e^{imx}$
 - Both l and m are in the range $-n \dots n$. **Meaning that?**
- In fact, for a given Fourier mode k :

$$\frac{d\hat{u}_k}{dt} = - \sum_{l=-\alpha}^{\alpha} \sum_{m=-\alpha}^{\alpha} i m [\hat{u}_l \hat{u}_m e^{i(l+m)x}] e^{-ikx}, \text{ for } -n \leq k \leq n$$

Computing individual terms in spectral space is computationally expensive, even in 1D

- α must be large enough to cover all possible combinations of l, m inside the range $-n, n$
- The number of points inside each triangular (eXcluded) region is:

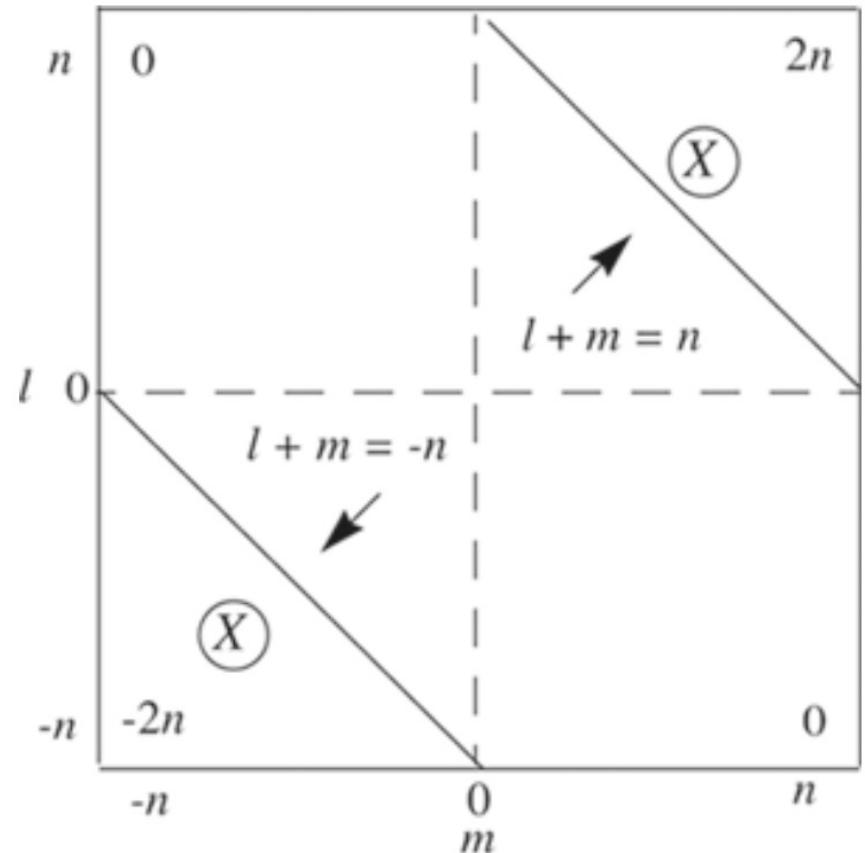
$$1 + 2 + 3 \dots + (n-1) = \frac{n(n-1)}{2}$$

- The number of points *retained* is therefore:

$$(2n+1)^2 - 2 \left[\frac{n(n-1)}{2} \right] = 3n^2 + 5n + 1$$

Which is the number of points that must be evaluated.

- This is $O(n^2)$ and it is very bad news, also considering that it originated with just a quadratic term. Computational cost will increase very rapidly. This is much worse in 2D.
- What can we do?



Combining grid and spectral methods

- What if we evaluated u and $\frac{\partial u}{\partial x}$ on a grid? If we had a sufficiently large grid, no aliasing!

$$\left(u \frac{\partial u}{\partial x} \right)_k = \frac{1}{M} \sum_{j=1}^M \left[u(x_j) \frac{\partial u}{\partial x}(x_j) e^{-ikx_j} \right], \quad -n \leq k \leq n$$

SPECTRALGRID

- Now use definitions to write terms as Fourier series (and benefit from exact derivatives)

$$\left(u \frac{\partial u}{\partial x} \right)_k = \frac{1}{M} \sum_{j=1}^M \left[\left(\sum_{l=-n}^n \hat{u}_l e^{ilx_j} \right) \left(\sum_{m=-n}^n i m \hat{u}_m e^{imx_j} \right) e^{-ikx_j} \right], \quad -n \leq k \leq n$$

SPECTRALGRID, using SPECTRAL METHOD

How many grid points do we need for the non-linear momentum advection?

- We have 3 Fourier modes, k, l, m , so zonal wave numbers up to $\pm 3n$. Therefore we need $3n+1$ complex coefficients.
- But $u \frac{\partial u}{\partial x}$ is real, so we only have $3n+1$ independent real numbers.
 - Therefore $M \geq 3n+1$
- 50% more expensive than the linear case, but not as bad as n^2
- Note that the extra 50% of information on the grid is thrown away when we transform back (5)

Our spectral model so far:

1. Initialise spectral coefficients \hat{u}_k for $-n \leq k \leq n$
1. Evaluate u and $\frac{\partial u}{\partial x}$ on a grid with M points.

The derivative is computed using the spectral method
2. Form the product $u \frac{\partial u}{\partial x}$ on the grid
3. Now transform back into wave-number space, so as to obtain coefficients:
$$\left(\widehat{u} \frac{\partial u}{\partial x} \right)_k$$
5. Predict (do the time step)
6. Go back to step 2.
$$\frac{d\hat{u}_k}{dt} = - \left(\widehat{u} \frac{\partial u}{\partial x} \right)_k$$

Spectral methods on the sphere

- First developed by Silberman (1954)

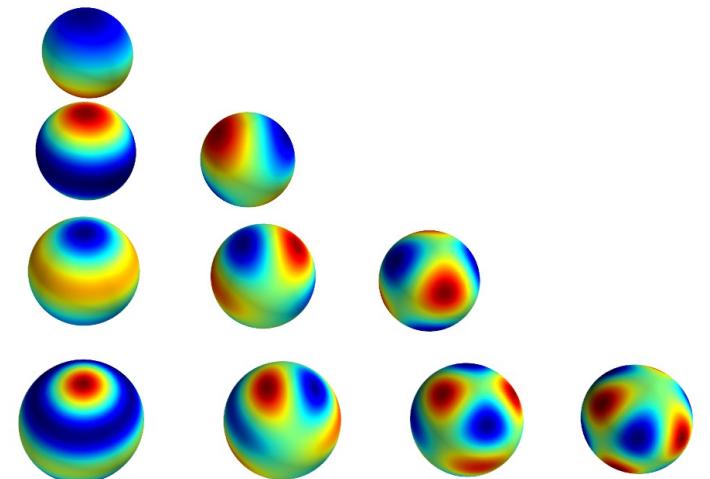
$$F(\lambda, \varphi) = \sum_{m=-\infty}^{\infty} \sum_{n=|m|}^{\infty} F_n^m Y_n^m(\lambda, \varphi),$$

λ is lon; φ is lat; m is the zonal wavenumber

where:

$$Y_n^m(\lambda, \varphi) = e^{im\lambda} P_n^m(\sin \varphi)$$

are spherical harmonics and $P_n^m(\sin \varphi)$ are the associated Legendre functions of the first kind



The spherical harmonics Y_n^m are eigenfunctions of the Laplacian on the sphere:

$$\nabla^2 Y_n^m = \frac{-n(n+1)}{a^2} Y_n^m \quad \text{where } a \text{ is the radius of the sphere}$$

In practice: spectral models on the sphere

An approximation to F can be made by using truncated sums:

$$\bar{F} = \sum_{m=-M}^M \sum_{n=|m|}^{N(m)} F_n^m Y_n^m \quad m \text{ is the zonal wavenumber}$$

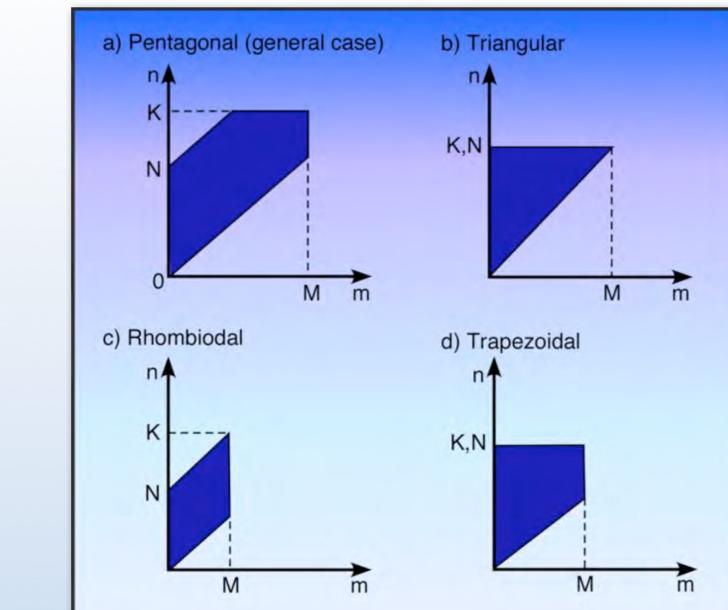
how to choose $N(m)$?

This is known as the problem of *truncation*.

Two well-known possibilities are:

rhomboidal truncation $N=M+|m|$ and
triangular truncation $N=M$

Different Truncation Types used in spectral AGCM



[McGuffie & Henderson-Sellers, A Climate Modelling Primer]

Spherical Harmonic Transform

- Similar to what we did with Fourier transforms, we can design a spherical harmonic transform, which consists in the combination of a Fourier transform and a Legendre transform.
- The Legendre transform is formulated using a method called *Gaussian quadrature*.
 - Say that we want to evaluate: $I \equiv \int_{-1}^1 f(x) dx$ numerically
 - If $f(x)$ is defined at a finite number of points x_i , then $I \cong \sum_{i=1}^N f(x_i) w_i$, w_i are weights
- Special case: $f(x)$ is a weighted sum of Legendre polynomials, as in a Legendre transform. We want the transform to be exact.
- Gauss showed that (44) gives the exact value of I if we choose the points x_i to be the roots of the highest Legendre polynomial used.
- **In practice, we choose (Gaussian) latitudes to be those roots.** Expensive (iterative), but we only need to perform this calculation once.

Spherical grids

- For either truncation, rhomboidal or triangular, choosing M fixes the expansion. That is then how we can say R15 or T106. The number of coefficients is:
 - Rhomboidal: $(M+1)^2 + M^2 + M$
 - Triangular: $(M+1)^2$
- So, the transform suitable for a spherical geometry needs many more points than the spectral transform. For triangular truncation:
 - Grid points around a latitude circle: $\geq 3M+1$
 - Grid points on a meridian: $\geq (3M+1)/2$
 - **TOTAL:** $\geq (3M+1)^2/2$ a factor of 2.25 between grid and spectral representations.
- Computing physics on the non-aliasing grid is quite wasteful, but unavoidable.
- Until recently we did not have fast Legendre transforms, but ECMWF have made major breakthroughs in this area. Without FLTs, operation count would be $O(N^3)$, while finite difference methods are $O(N^2)$.

Summary of spectral methods

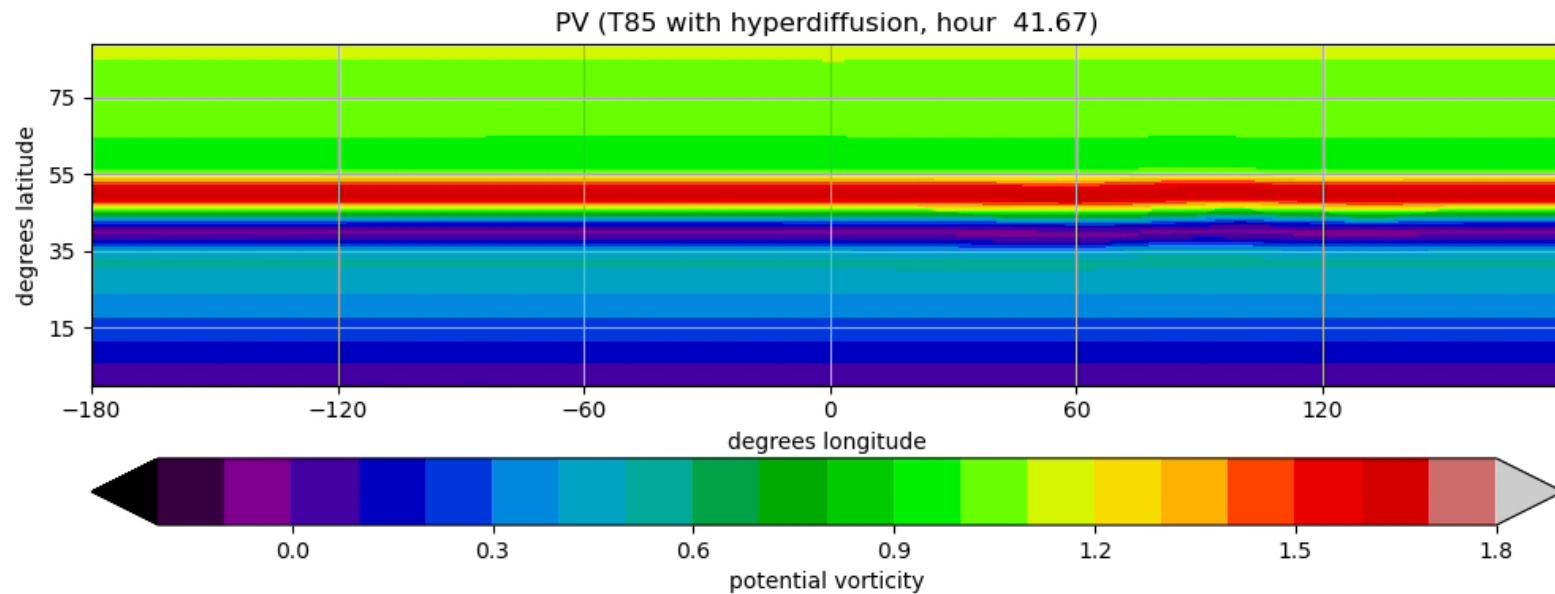
Strengths

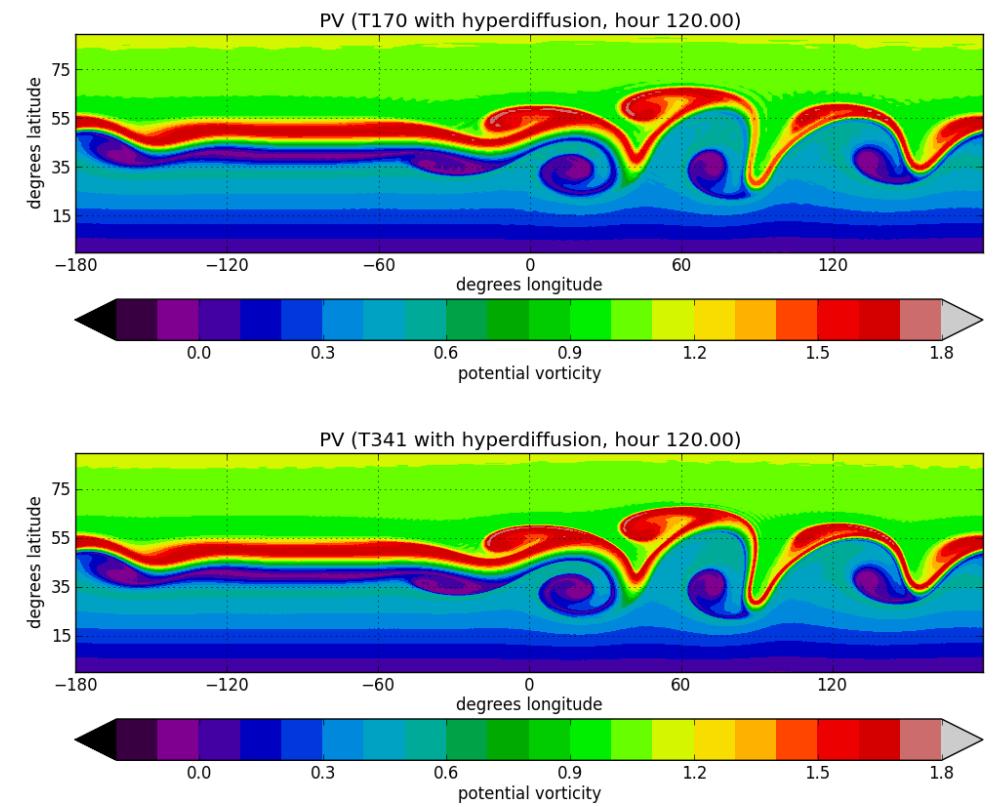
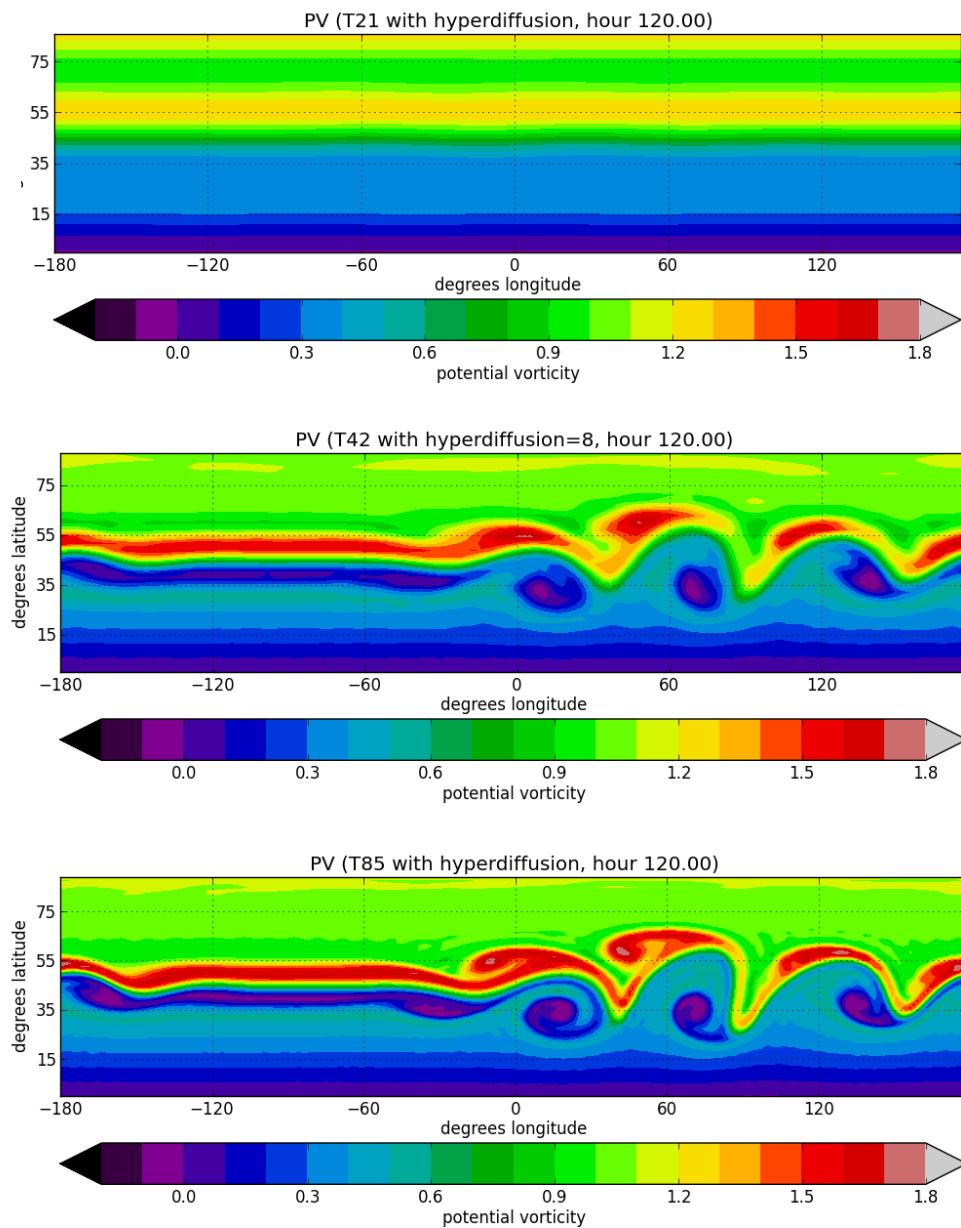
- Eliminate pole problem, especially with triangular truncation
- Provides exact phase speed for linear waves and advection by constant current
- **Converges very rapidly and gives good results even with just few modes**
- Semi-implicit time stepping is implemented very easily

Weaknesses

- Do not conserve anything
- Failure to conserve requires damping, else BOOM
- “Bumpy” oceans
- Truncation in transform may mean that physics do not always have the intended effects
- Moisture advection problematic
- Until FLT, not practical at high-resolution

Galewsky's barotropic instability problem with a global spectral model (T85)





The Galewski problem on truncations T21 to T341 at time 120 hours.

T42 already captures the main characteristics of the flow.

Results from a global, spectral SWE model that was written in Python and can be run on a laptop

A spectral SWE model

1) Setup grid and initial conditions

```
# setup up spherical harmonic instance, set lats/lons of grid
x = Spharmt(nlons, nlats, ntrunc, rsphere, gridtype="gaussian")
# x = Spharmt(nlons, nlats, ntrunc, rsphere, gridtype="regular")
lons, lats = np.meshgrid(x.lons, x.lats)
f = 2.*omega*np.sin(lats) # coriolis

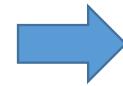
# zonal jet.
vg = np.zeros((nlats, nlons), np.float)
u1 = (umax/en)*np.exp(1./((x.lats-phi0)*(x.lats-phi1)))
ug = np.zeros((nlats), np.float)
ug = np.where(np.logical_and(x.lats < phi1, x.lats > phi0), u1, ug)
ug.shape = (nlats, 1)
ug = ug*np.ones((nlats, nlons), dtype=np.float) # broadcast to shape (nlats, nlons)
# height perturbation.
hbump = hamp*np.cos(lats)*np.exp(-((lons-np.pi)/alpha)**2)*np.exp(-(phi2-lats)**2/beta)

# initial vorticity, divergence in spectral space
vrtspec, divspec = x.getvrtdivspec(ug, vg)
vrtg = x.spectogrd(vrtspec)
divg = x.spectogrd(divspec)

# create hyperdiffusion factor
hyperdiff_fact = np.exp((-dt/efold)*(x.lap/x.lap[-1])**2*(ndiss/2))

# solve nonlinear balance eqn to get initial zonal geopotential,
# add localized bump (not balanced).
vrtg = x.spectogrd(vrtspec)
tmpg1 = ug*(vrtg+f)
tmpg2 = vg*(vrtg+f)
tmpspec1, tmpspec2 = x.getvrtdivspec(tmpg1, tmpg2)
tmpspec2 = x.grdtospec(0.5*(ug**2+vg**2))
phispec = x.invlap*tmpspec1 - tmpspec2
phig = grav*(hbar + hbump) + x.spectogrd(phispec)
phispec = x.grdtospec(phig)

# initialize spectral tendency arrays
ddivdtspec = np.zeros(vrtspec.shape+(3,), np.complex)
dvrtdtspec = np.zeros(vrtspec.shape+(3,), np.complex)
dphidtspec = np.zeros(vrtspec.shape+(3,), np.complex)
nnew = 0
nnow = 1
nold = 2
```



2) Time integration loop

```
# time loop.
time1 = time.time()
for ncycle in range(itmax+1):
    t = ncycle*dt
    # get vort, u, v, phi on grid
    vrtg = x.spectogrd(vrtspec)
    ug, vg = x.getuv(vrtspec, divspec)
    phig = x.spectogrd(phispec)
    print("t=%6.2f hours: min/max %6.2f, %6.2f" % (t/3600., vg.min(), vg.max()))

    # compute tendencies.
    tmpg1 = ug*(vrtg+f)
    tmpg2 = vg*(vrtg+f)
    ddivdtspec[:, nnew], dvrtdtspec[:, nnew] = x.getvrtdivspec(tmpg1, tmpg2)
    dvrtdtspec[:, nnew] *= -1

    tmpg = x.spectogrd(ddivdtspec[:, nnew])
    tmpg1 = ug*phig
    tmpg2 = vg*phig
    tmpspec, dphidtspec[:, nnew] = x.getvrtdivspec(tmpg1, tmpg2)
    dphidtspec[:, nnew] *= -1

    tmpspec = x.grdtospec(phig+0.5*(ug**2+vg**2))
    ddivdtspec[:, nnew] += -x.lap*tmpspec

    # update vort, div, phiv with third-order adams-bashforth.
    # forward euler, then 2nd-order adams-bashforth time steps to start.
    if ncycle == 0:
        dvrtdtspec[:, nnow] = dvrtdtspec[:, nnew]
        dvrtdtspec[:, nold] = dvrtdtspec[:, nnew]
        ddivdtspec[:, nnow] = ddivdtspec[:, nnew]
        ddivdtspec[:, nold] = ddivdtspec[:, nnew]
        dphidtspec[:, nnow] = dphidtspec[:, nnew]
        dphidtspec[:, nold] = dphidtspec[:, nnew]
    elif ncycle == 1:
        dvrtdtspec[:, nold] = dvrtdtspec[:, nnew]
        ddivdtspec[:, nold] = ddivdtspec[:, nnew]
        dphidtspec[:, nold] = dphidtspec[:, nnew]

        vrtspec += dt*(
            (23./12.)*dvrtdtspec[:, nnew] - (16./12.)*dvrtdtspec[:, nnow]
            + (5./12.)*dvrtdtspec[:, nold])
        divspec += dt*(
            (23./12.)*ddivdtspec[:, nnew] - (16./12.)*ddivdtspec[:, nnow]
            + (5./12.)*ddivdtspec[:, nold])
        phispec += dt*(
            (23./12.)*dphidtspec[:, nnew] - (16./12.)*dphidtspec[:, nnow]
            + (5./12.)*dphidtspec[:, nold])

    # implicit hyperdiffusion for vort and div.
    vrtspec *= hyperdiff_fact
    divspec *= hyperdiff_fact
    # switch indices, do next time step.
    nsav1 = nnew
    nsav2 = nnow
    nnew = nold
    nnow = nsav1
    nold = nsav2
```

Bring prognostics to grid

Tendencies on grid

Tendencies
in spectral
space

Adams-Bashfort
Time step

Physics in
spectral space

Update

Equivalent grid resolution of spectral models

We will see later in the lecture that, in 1D, we need at least $2N+1$ grid points to represent the waves used in a truncation with N wavenumbers.

The real resolution needs 1D lie between $2N+1$ and N^2 , as also seen in the paper by Renee Laprise (1992).

However, what is the real, effective resolution of (spectral) models? We have a paper under review (Klaver et al. 2019) to answer that question.

Laprise: 4 ways to compute the resolution of a spectral model. Let us take T31 truncation

1: average distance between latitudes on Gaussian grid $L1 = 425\text{km}$

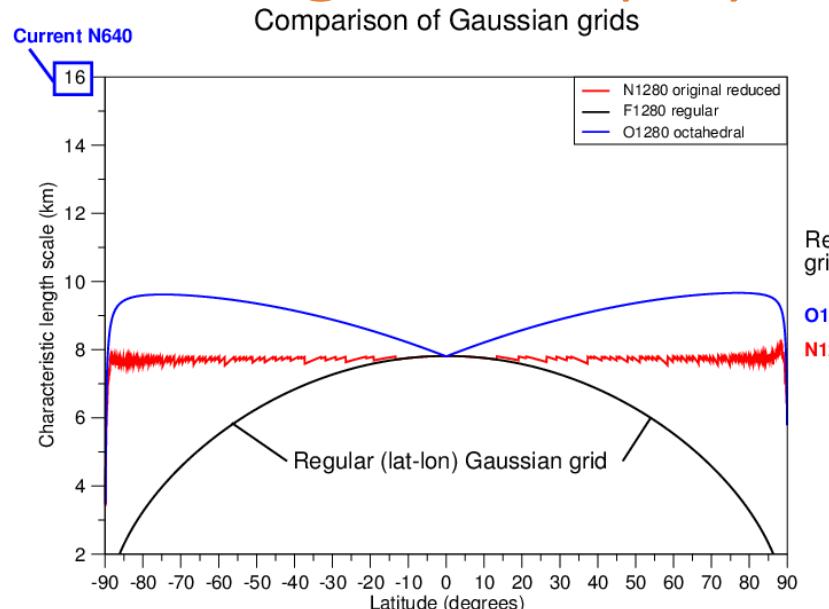
2: half the wavelength of the shortest resolved zonal wave at the equator: $L2 = 650\text{km}$

3: Number of real variables per unit area: $L3 = 725\text{km}$

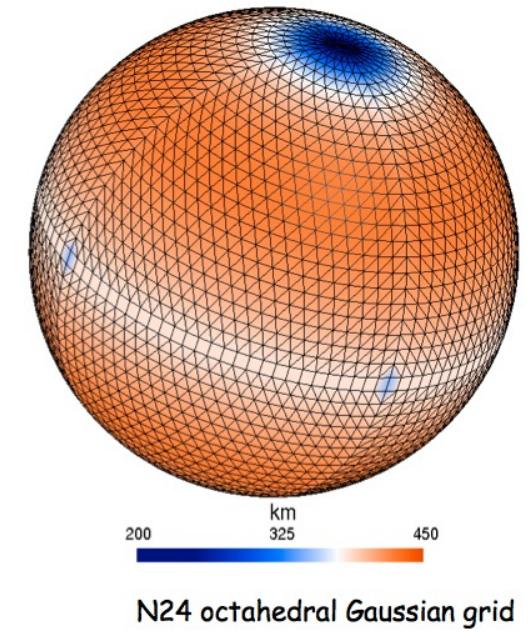
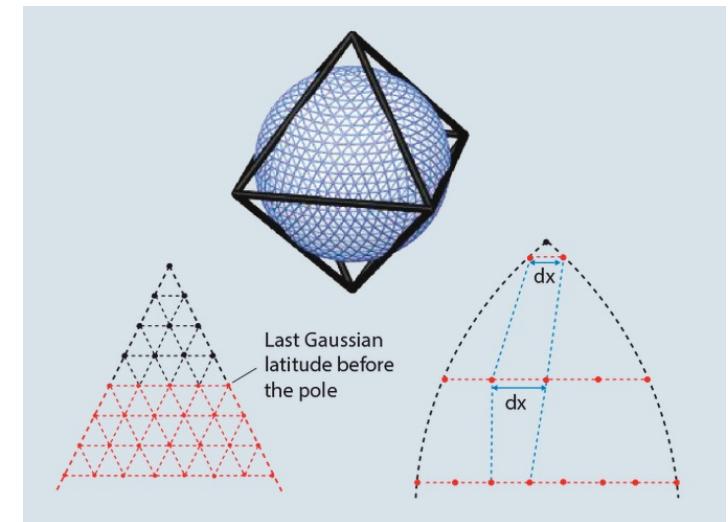
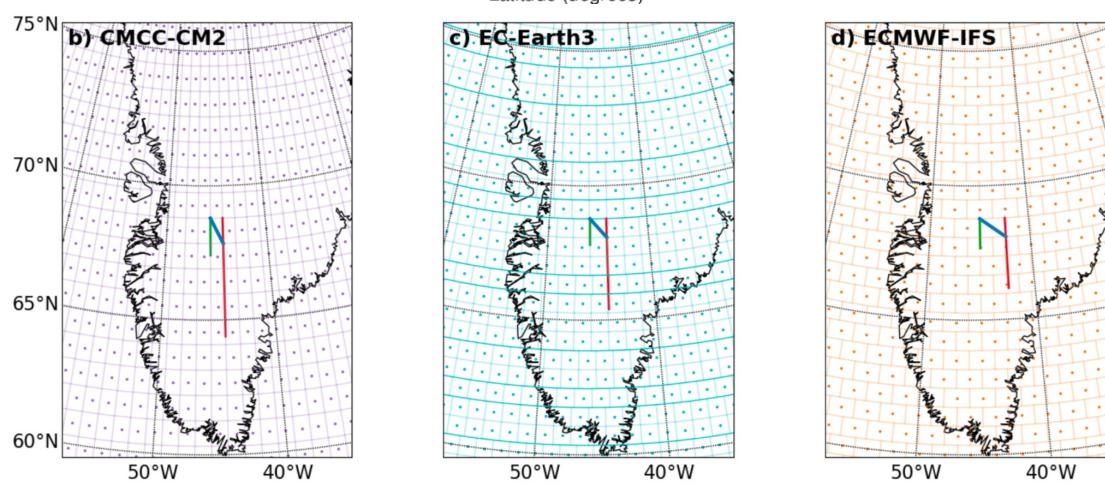
4: equivalent total wave number associated with the Laplacian operator (for highest mode): $L4 = 900\text{km}$

For reference, in lat/lon models, $N31 = 425\text{km}$

Associated grids in physical space

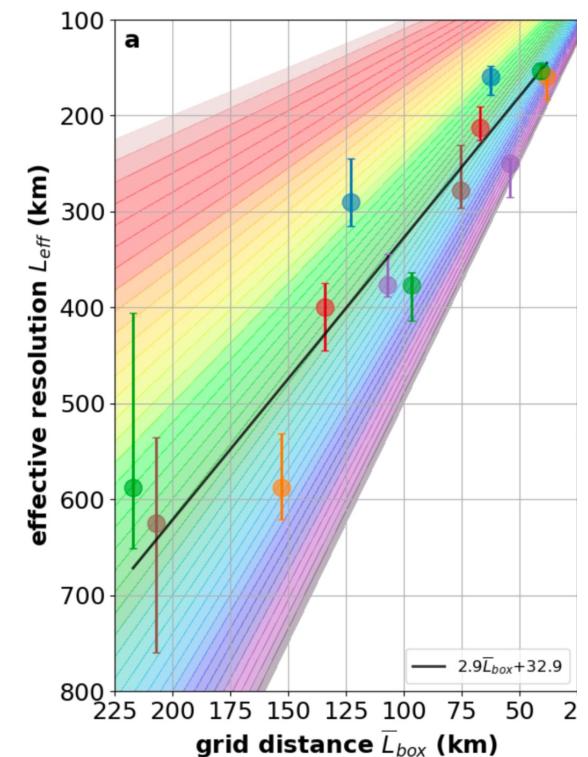


Reduced grids:
O1280
N1280

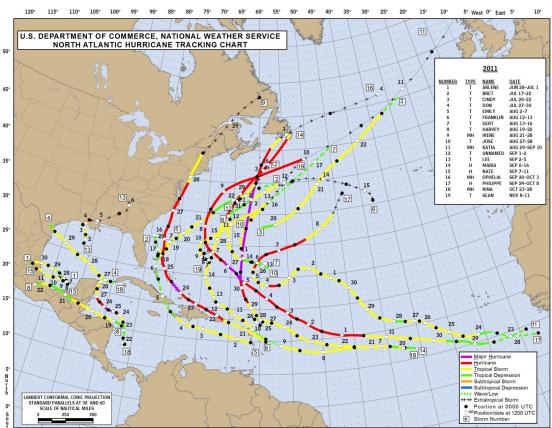
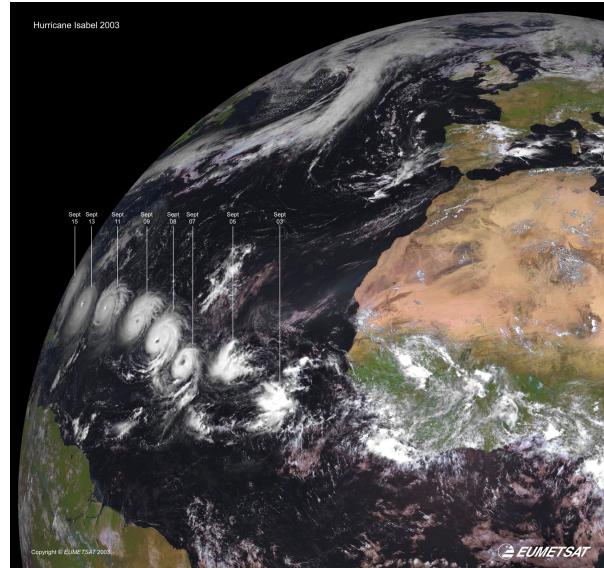


Effective resolution of PRIMAVERA models

Model-Version	Grid Type	Native Grid	\bar{L}_{box} (km)	Resolution L_{eff} (km)	L_{eff}/\bar{L}_{box}
HadGEM3-GC31-LM	grid point	145 x 192	217	525-770 (588)	2.4-3.5 (2.7)
HadGEM3-GC31-MM	grid point	325 x 432	96.7	340-390 (377)	3.5-4.1 (3.9)
HadGEM3-GC31-HM	grid point	769 x 1024	40.8	145-160 (153)	3.6-3.9 (3.7)
CMCC-CM2-HR4	grid point	192 x 288	153	555-645 (588)	3.6-4.2 (3.8)
CMCC-CM2-VHR4	grid point	768 x 1152	38.2	135-165 (159)	3.5-4.4 (4.2)
ECMWF-IFS-LR	spectral	reduced TCO255	123	265-335 (290)	2.2-2.7 (2.4)
ECMWF-IFS-HR	spectral	reduced TCO511	62.6	140-170 (159)	2.2-2.7 (2.5)
EC-Earth3	spectral	reduced TL255	107	365-410 (377)	4.0-4.7 (4.6)
EC-Earth3-HR	spectral	reduced TL511	54.2	215-255 (250)	3.4-3.8 (3.5)
MPIESM-1-2-HR	spectral	TQ127	134	355-425 (400)	2.7-3.2 (3.0)
MPIESM-1-2-XR	spectral	TQ255	66.9	200-235 (213)	3.0-3.5 (3.2)
CNRM-CM6-0	spectral	reduced TL127	207	490-715 (625)	2.4-3.5 (3.0)
CNRM-CM6-0-HR	spectral	reduced TL359	75.3	260-325 (278)	3.4-4.3 (3.7)



Finding the storms

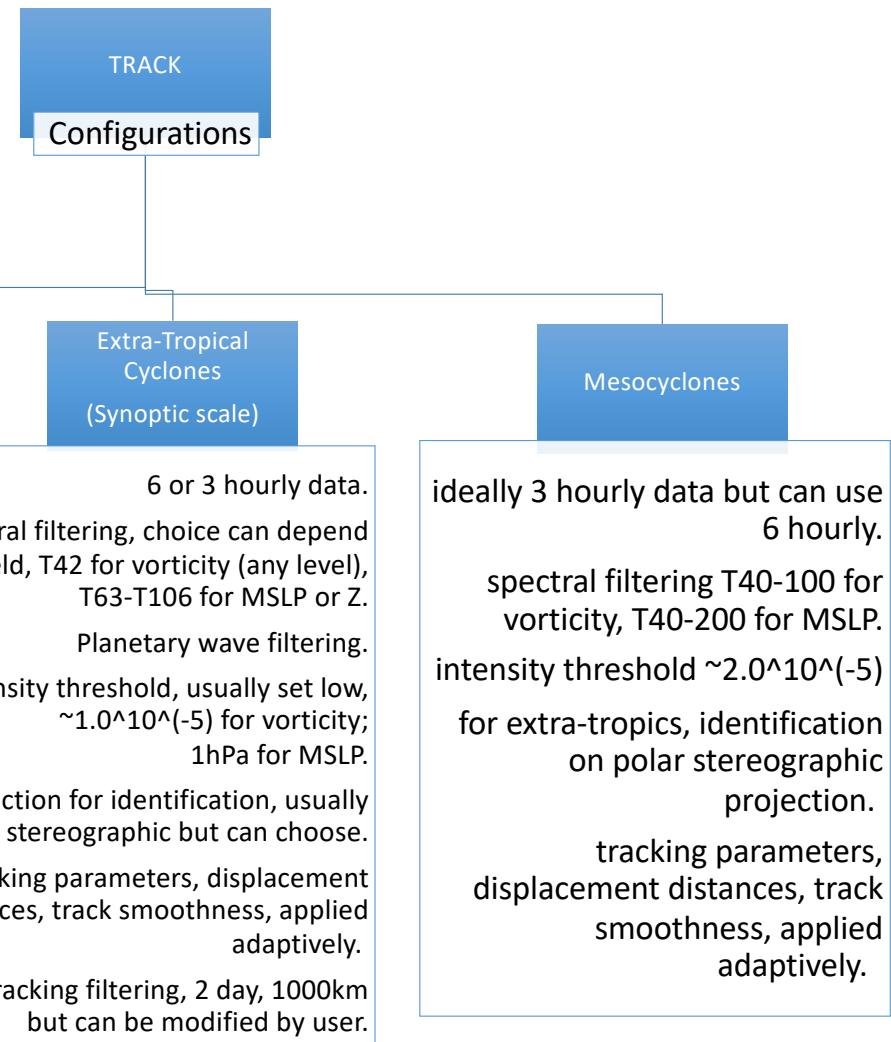


6 hourly data.
spectral filtering, T42 for single level vorticity; T63 for vertically averaged vorticity.

Planetary wave filtering.
intensity threshold, usually set low, $\sim 0.5 \times 10^{-5}$.

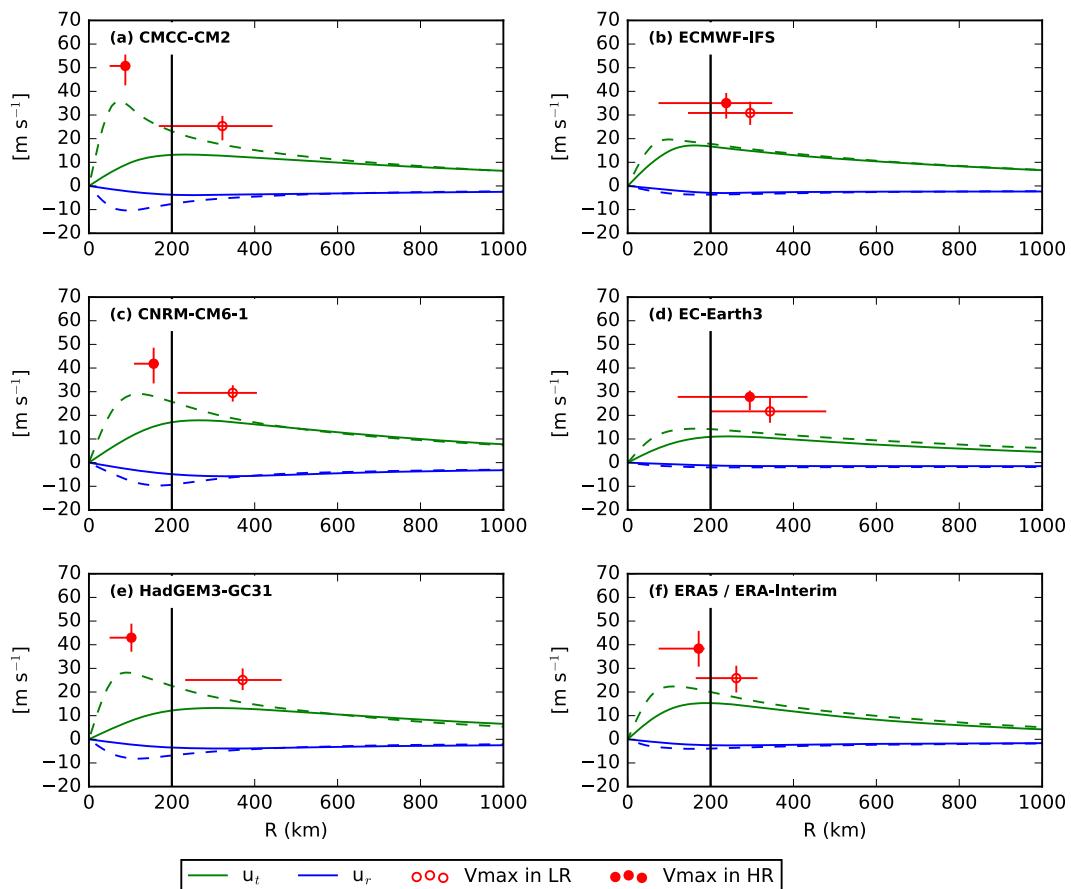
cylindrical projection (60S, 60N).
tracking parameters, displacement distances, track smoothness, applied adaptively.

post tracking filtering, 2 days.
warm core test, add vorticity at T63 for multiple levels up to 200hPa.



Extra slides

Simulation of the primary and secondary circulations of hurricanes



PRIMAVERA GCMs:
composites of Tropical
Cyclone structures at high
and low resolution

The inner core region is better simulated at high resolution. This is clearer in finite difference and finite volume models (left column, top and bottom) than it is in spectral models (right column and left-centre, CNRM)