# Simulating seismic wave propagation with sw4

CIG tutorial

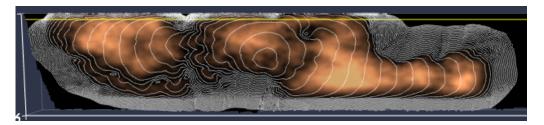
N. Anders Petersson



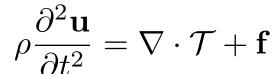


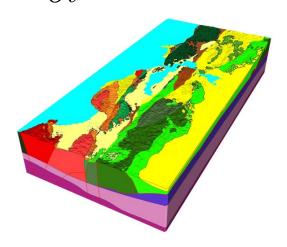
#### SW4 simulates regional earthquake ground motions on parallel computers

- Mathematical model of the Earth:
  - Navier's eqn of 3-D linear isotropic solid dynamics
  - 4<sup>th</sup> order summation-by-parts finite differences
  - 4<sup>th</sup> order time explicit stepping
- Geophysical model
  - Isotropic heterogeneous elastic & visco-elastic material
  - Topography
- Kinematic source rupture model



- MPI-based distributed memory programming model
  - C++ with Fortran numerical kernels
- SW4 is open source, GNU GPL-v2

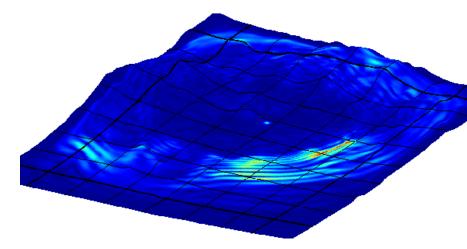






## The SW4 and WPP codes solve the time dependent 3D seismic wave equation

- WPP is 2<sup>nd</sup> order accurate and SW4 is 4<sup>th</sup> order
- Summation by parts (SBP) finite difference methods
  - Fundamentally different from FDTD
  - Energy stability
  - Higher order accuracy
- Curvilinear coordinates
  - Realistic topography
  - Absorbing far-field layers
- Visco-elastic modeling
- Point forces and moment tensor sources
  - Grid independent location
- Anisotropic materials
- Local mesh refinement



### Summation by parts discretization mimic integration by parts for finite differences

• Elastic wave equation:

$$\rho \frac{\partial^2 \mathbf{u}}{\partial t^2} = \mathbf{L}\mathbf{u} + \mathbf{f}(\mathbf{x}, t)$$
$$\mathbf{L}\mathbf{u} := \nabla \cdot \mathcal{T}$$

Discretized in space:

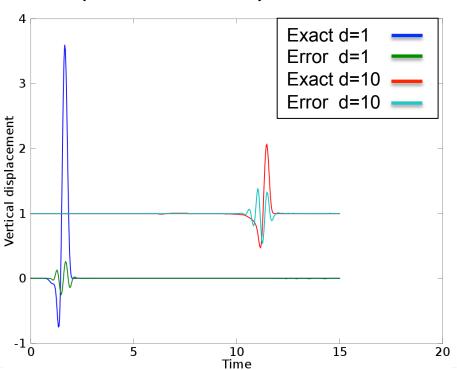
$$\rho \frac{d^2 \mathbf{u}}{dt^2} = \mathbf{L}_h \mathbf{u} + \mathbf{f}(\mathbf{x}, t)$$

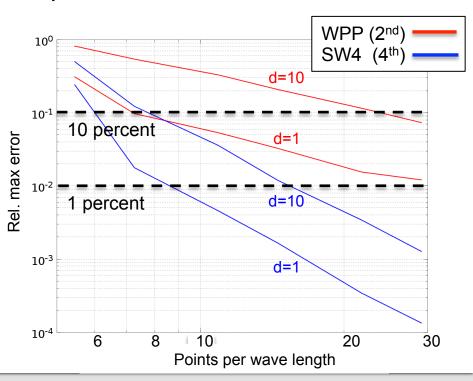
- Symmetry and negative definite properties of 'L' preserved by summation by parts discretization 'L<sub>h</sub>'
- Numerical scheme is energy stable



A 4<sup>th</sup> order method is more efficient since it needs fewer grid points per wavelength

- Lamb's problem
- Solution at distance d=1 and d=10 [Mooney (1974)]
- Error: P<sup>-2</sup> versus P<sup>-4</sup> (P = grid points per wave length)
- Improved efficiency with 4<sup>th</sup> order accuracy





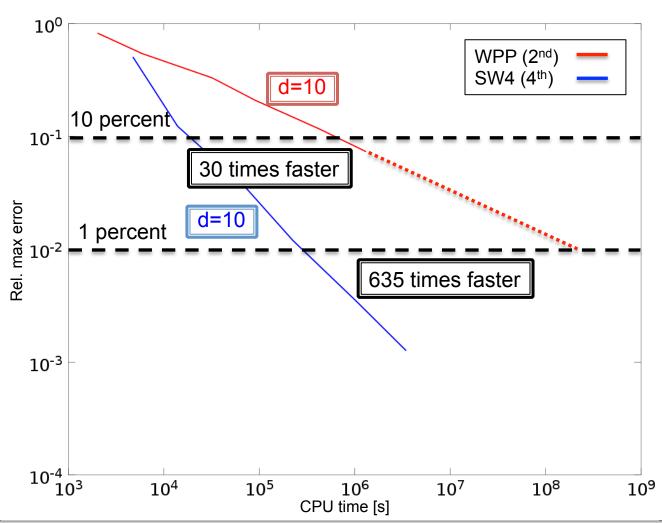




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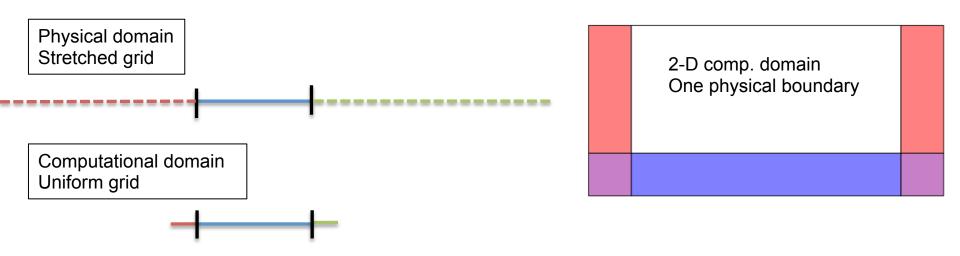
# The 4<sup>th</sup> order method is much faster at larger distances and gives smaller errors

#### Error at distance d=10 from source

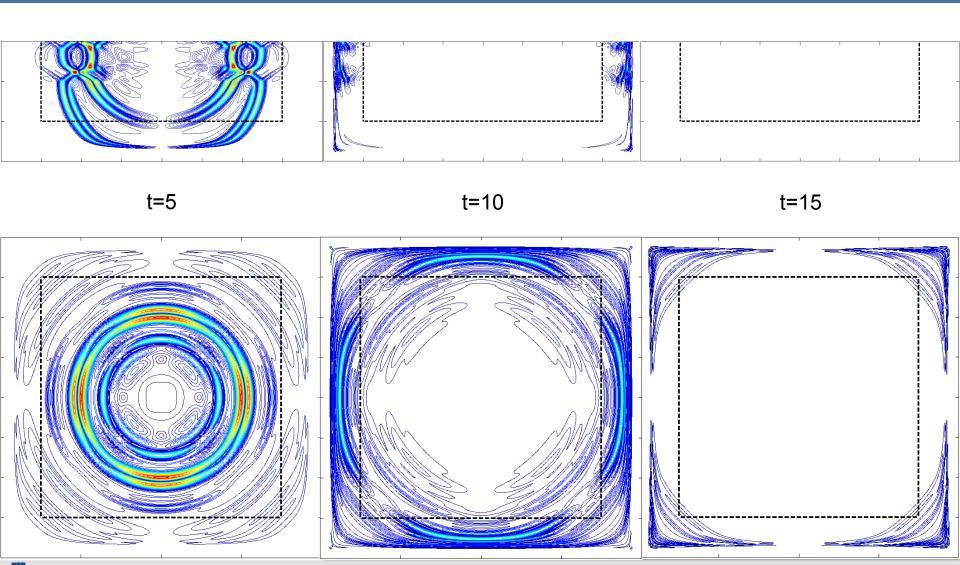


### SW4 uses super-grid layers to suppress artificial reflections from far-field boundaries

- Perfectly matched layers (PML) can be unstable
- Super-grid shrinks very large physical domain
  - Slow down and compress waves in layers
  - Damp unresolved modes with high order dissipation
- Always stable
  - Accurate if layer is sufficiently wide



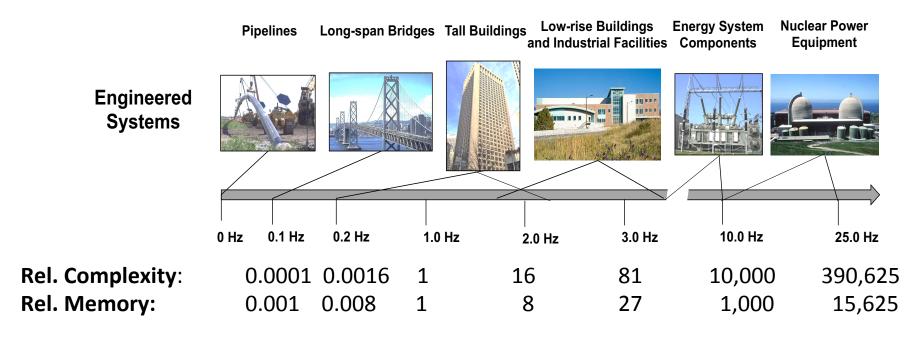
# Super-grid test: 3-D elastic wave equation with layered material and a free surface





# The computational requirements scale as (frequency)<sup>4</sup>

- Smallest wave length  $\lambda \approx V_s/f_{max}$
- Grid size:  $h = \lambda/P$ ; Points per wavelength: P
- Double frequency -> 8 x grid points, 16 x CPU



#### Hands-on sw4 tutorial

#### How to run sw4

Smaller & larger jobs

#### How to set up the simulation

- Sources
- How do I pick the grid size?
- Topography
- Material model, visco-elastic modeling
- Output options

#### Practical suggestions

- Start coarse, check model, estimate resources, refine
- Use the parallel file system
- Initial analysis before moving large files

#### Lots of info in the user's guide

— VM: ~/Documents/SW4-UsersGuide.pdf

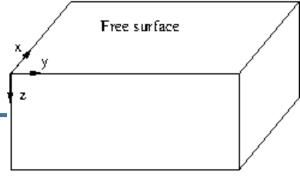
#### Building and running sw4

- sw4 can be built on Linux/OSX/Unix machines
  - ~/Documents/SW4-Installation.pdf
- Command file specifies the simulation
  - mpirun -np 2 sw4 my-input.in
- Syntax for running MPI-jobs varies:
  - Macs with openmpi: openmpirun –np 4 sw4 input.in
  - VM-ware mpirun: mpirun –np 8 sw4 input.in
  - Livermore computing (LC): srun -n4096 sw4 my-big-job.in
- Running on supercomputers (details may vary)
  - Smaller, shorter jobs on interactive debug partition. Start job from command prompt. No CPU-bank needed
  - Larger jobs use a batch system. Job described by a script. Usually needs a CPU-bank
  - For faster I/O, use a parallel file system if available

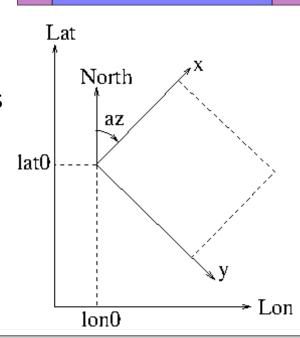


# **sw4** uses SI-units (MKS), angles in degrees

- Box-shaped computational domain
  - Right-handed system, z positive downwards
- Top surface optionally follows topography
- Free surface boundary condition on top surface
- Super-grid absorbing layers on all other sides
  - SG layers included in computational domain
  - Default thickness: 30 grid points
  - Smallest Cartesian grid: 60x60x30
- Locations in Cartesian or geographic coordinates
- Geographic coordinates are mapped:
  - Default: spheroidal projection
  - Proj.4 library for better accuracy
- Rotate grid with azimuth angle (az)
  - az=0: x-axis = North, y-axis = East



All sources and receivers in interior



#### 1st exercise: solve Lamb's problem

- Start VM virtualBox & import the CIG16-TutorialVM appliance
- Log on as sw4\_user (No password)
- cd Desktop
- Simulation described in input file: seismicl.in fileio path=seis-out grid nx=151 x=10e3 y=10e3 z=5e3 time t=5.0 # supergrid thickness for h = 10000/150 = 1000/15, 30\*h = 2000, 60\*h = 4000 supergrid gp=60 block vp=1.7320508076e+03 vs=1000 rho=1500 source type=C6SmoothBump x=5e3 y=5e3 z=0 fz=1e13 freq=1 t0=0 # Time history of solution on the surface rec x=5e3 y=6e3 z=0 file=v1s sacformat=0 usgsformat=1
- mpirun -np 1 sw4 seismic1.in
- (1 min, 51 sec runtime on my Mac laptop)

#### We use octave to view the results

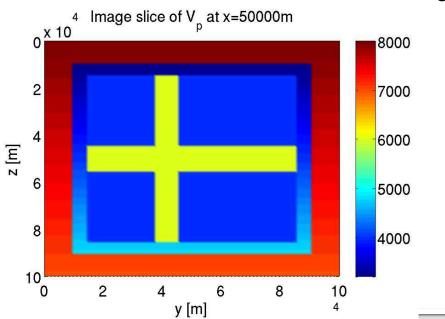
- octave is open source software (similar to Matlab)
- Open octave from terminal window (in VM):
  - >octave
- Expand octave search path (put command in ~/.octaverc file)
- octave:1> addpath("/home/sw4\_user/Desktop/tools")
- octave:2>cd Desktop/seis-out
- octave:3>ls
- octave:4>plotusgs("v1s.txt")

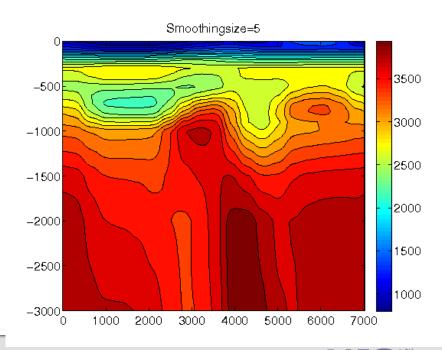
# The material model ( $\rho$ , $C_p$ , $C_s$ ) must be defined for all points in the grid

- One or several material commands can be combined
  - block, rfile, efile, pfile, ifile, ...
- The order matters (later commands take precedence)

Visco-elastic modeling triggered by attenuation command:

Also need to specify Q<sub>p</sub> and Q<sub>s</sub>



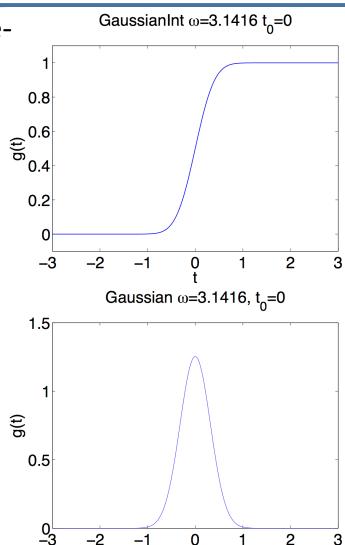


### **sw4** provides several options for specifying sources and time-functions

- Any number of point forces or moment tensor sources
  - Forces defined by 3-component vector
  - Moment tensor (symmetric) defined by 6 components M<sub>ii</sub>, or
  - Seismic moment (M<sub>0</sub>) and strike, dip, rake angles
- Many pre-defined time functions (Gaussian, Ricker, Liu, ...)
  - Frequency parameter and start/center time
- Dirac delta-distribution: trigger all frequencies on mesh
  - Discrete Green's functions. Motion must be filtered for accuracy
- User defined discrete time function (interpolated by spline)
- Complex ruptures can be specified by using many point sources or an SRF file

# There are several ways of computing displacements, velocities, or acceleration

- Displacements correspond to source timefunctions that tends to a constant (e.g. GaussianInt)
  - Displacements obtained directly
  - Velocities by differentiating once
  - Acceleration by differentiating twice
- Velocities correspond to time-functions that tend to zero but have non-zero integral (e.g. Gaussian)
  - Velocities obtained directly
  - Displacements by integrating once
  - Acceleration by differentiating once



### The highest frequency and the lowest wave speed determine the grid size

- Shortest wave length L<sub>min</sub>=min C<sub>S</sub>/f<sub>max</sub>
- Grid points per shortest wave length  $P_{min} = L_{min}/h$
- Good accuracy when P<sub>min</sub> > 6 to 10 (depending on distance)
- 3 approaches:
  - Tune freq in source time function to not trigger unresolved waves
  - Calculate  $f_{max} = min C_S/(P_{min} h)$ , filter out unresolved frequencies afterwards (sac)
  - Use prefilter command to pre-process all sources (not discussed today)
- Relation between  $f_{max}$  and frequency parameter (freq) is different for each time function (see UG 4.4)
- octave scripts for plotting time function and Fourier transform: fcnplot, ftfcnplot



# 2<sup>nd</sup> exercise: solve a layer over half-space problem. Input file: loh1-slow.in

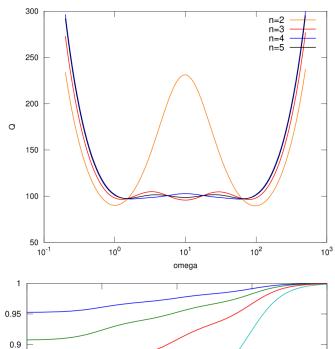
- The order of the block commands is important block vp=6000 vs=3464 rho=2700 block vp=4000 vs=2000 rho=2600 z2=1000
- Gaussian time function gives the velocities directly source x=15000 y=15000 z=2000 mxy=1e18 t0=1.2 freq=5.0 type=Gaussian
- Time histories recorded @ the closest grid point rec x=15600 y=15800 z=0 file=sta01 usgsformat=1 rec x=21000 y=23000 z=0 file=sta10 usgsformat=1
- Material images don't change, recorded initially image mode=s y=15e3 file=ver cycle=0
- Solution images can be recorded at fixed time intervals image mode=mag z=0 file=surf timeInterval=0.5

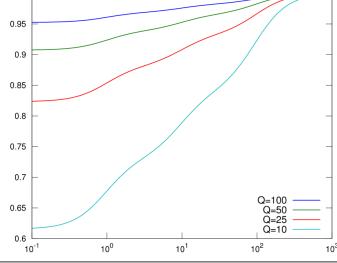
#### We can use octave to view image files

- octave>cd Desktop/loh1-slow
- octave>ls
- octave>help plotimage
- octave>plotimage("surf.cycle=170.z=0.mag.sw4img")
- octave>help imageinfo
- octave>imageinfo("surf.cycle=170.z=0.mag.sw4img",10,1)

### Standard linear solid (SLS) mechanisms are used to model anelastic attenuation

- The quality factors ( $Q_p$  and  $Q_s$ ) are approximately constant in frequency band [0.01  $f_{max}$ ,  $f_{max}$ ]
  - User must specify upper frequency: f<sub>max</sub>
  - More mechanisms: less variation in Q, but more calculations. Default nmech=3
- The visco-elastic material is dispersive: phase velocity depends on frequency
  - User must specify phasefreq parameter:
     frequency where C<sub>P</sub> and C<sub>S</sub> are accurate.





### 3<sup>rd</sup> exercise: solve an anelastic layer over halfspace problem. Input file: loh3-slow.in

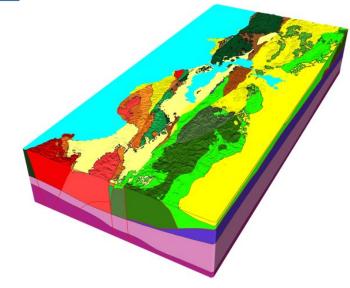
Only differences from loh1-slow.in

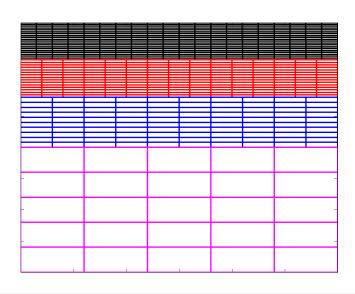
```
# different output directory
fileio path=loh3-slow
# must specify Qp and Qs
block vp=6000 vs=3464 rho=2700 Qs=69.3 Qp=155.9
block vp=4000 vs=2000 rho=2600 z2=1000 Qs=40 Qp=120
# enable attenuation modeling
attenuation phasefreq=2.5 nmech=3 maxfreq=15
```

About 3x CPU time and 2.7x memory

### Large complex material models can be described in an rfile model

- Variable resolution data structure
- 7.6 Gbyte file for regional SF bay area
- Vulcan (5 PFlops): 1 Gbyte memory per core
- Parallel I/O routines on parallel file systems





2897 x 1401 x 74 (6 Gb)

1449 x 701 x 57 (1.2 Gb)

725 x 351 x 33 (0.17 Gb)

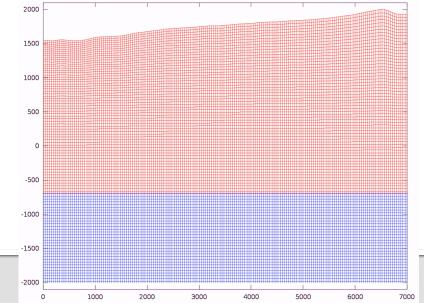
363 x 176 x 194 (0.25 Gb)

#### Realistic topography is included in an rfile

- SW4 constructs a curvilinear grid on startup
- z=0 is mean sea level
- User must pick bottom z-coordinate for curvilinear grid: zmax
- Rule of thumb: zmax = twice the variation below the lowest elevation
  - 1500 < elevation < 2500 m (positive above sea level)</li>

= zmax >= - (1500 – 2\*(2500-1500)) = +500 (positive below sea

level)



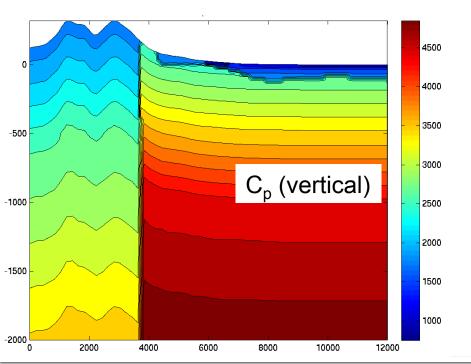
# 4<sup>th</sup> exercise: Simulate earthquake near Berkeley. Input file: berkeley-small.in

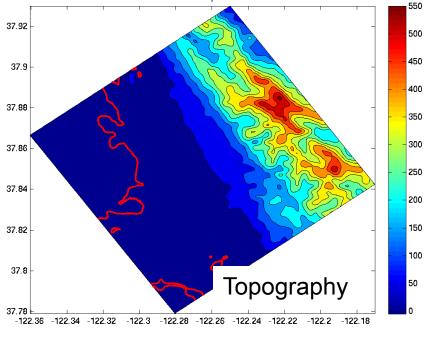
 Proj.4 commands in grid command (az must agree with rfile header)

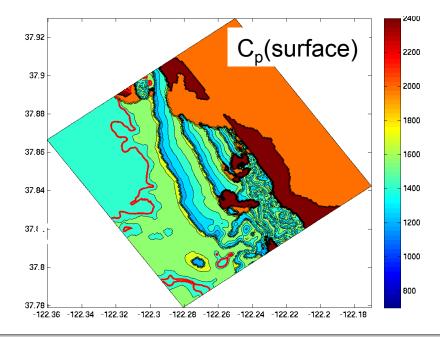
```
grid x=12e3 y=12e3 z=5e3 nx=151 lat=37.93 lon=-122.25 az=143.638 proj=tmerc datum=NAD83 lon p=-123.0 lat p=35.0 scale=0.9996
```

- Topography command topography input=rfile zmax=2e3 order=3 file=berkeley-r.rfile
- Material model rfile filename=berkeley-r.rfile
- Runtime 9 min, 50 sec

### The case berkeley-small.in 37.92 uses a small rfile model



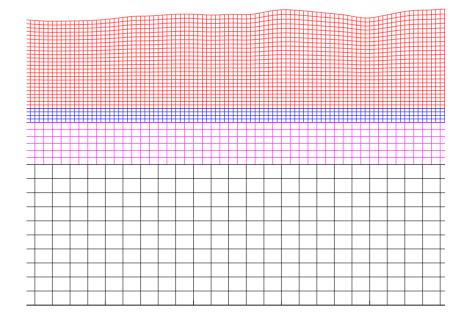






### Mesh refinement can be used to save computational resources

- Keep number of grid points per wave length ~ constant
- Saves memory
- Allows larger time step
- Currently only elastic materials



### 5<sup>th</sup> exercise: Try mesh refinement. Input file meshref.in

- The grid command specifies the coarsest grid size grid h=350 x=30000 y=30000 z=17000
- User specifies z-level with the refinement command refinement zmax=2.5e3
- Each grid patch must have at least 12 grid points in the vertical direction.
- Runtime ~5 min
- Same problem w/o mesh refinement in nomeshref.in

#### **General guidelines**

- Check setup on a coarse grid
- Check a few images of the material model
- Check receivers on a map: gmt
- Extrapolate computational resources from coarse run
- Doubling frequency = Halving the grid size = doubling # grid points / dimension
- Doubling # grid points/dimension: 8x grid points, 2x time steps
  - On same number of cores: 8x memory, 16x CPU time
  - On 8x more cores: same memory, 2x CPU time (weak scaling)
  - On 16x cores: ½ x memory, about the same CPU time
- SW4 source distribution includes examples and octave scripts

#### What if it doesn't work?

- Check command syntax and rules in SW4 user's guide
- Update SW4 source code from github
- Ask questions on CIG-SEISMO mailing list

cig-seismo@geodynamics.org

Discuss problems in the github issue tracker

https://github.com/geodynamics/sw4/issues

Help provided as time permits

