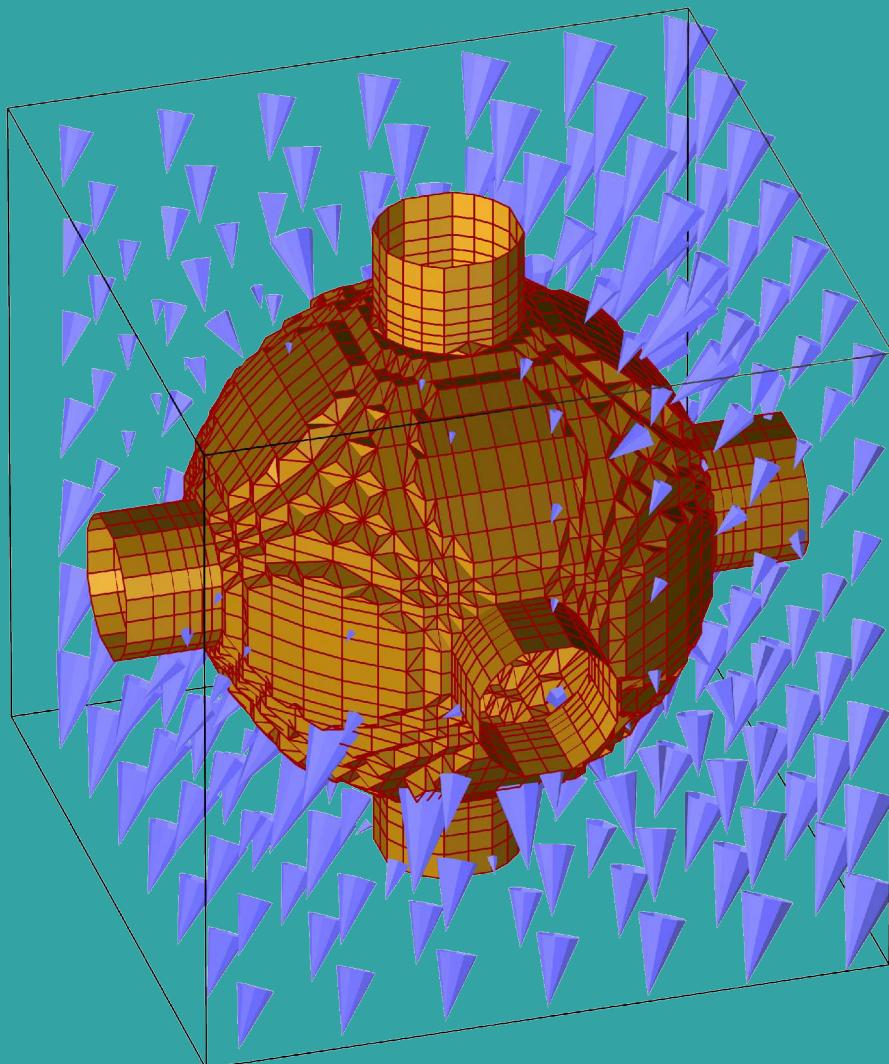
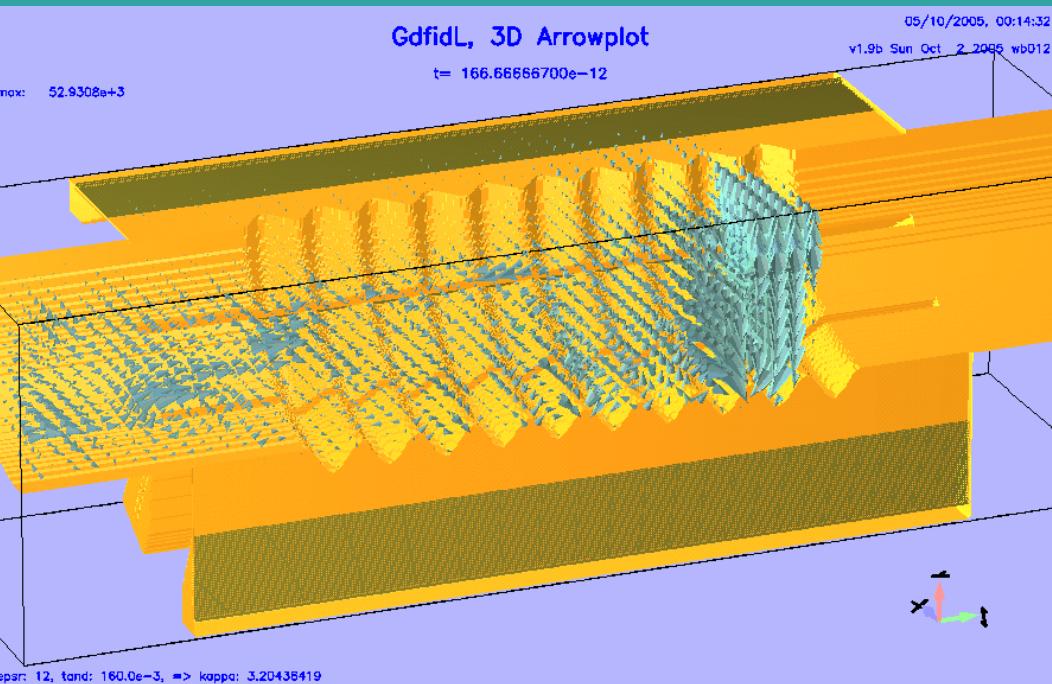


# The GdfidL Electromagnetic Field Simulator



Warner Bruns

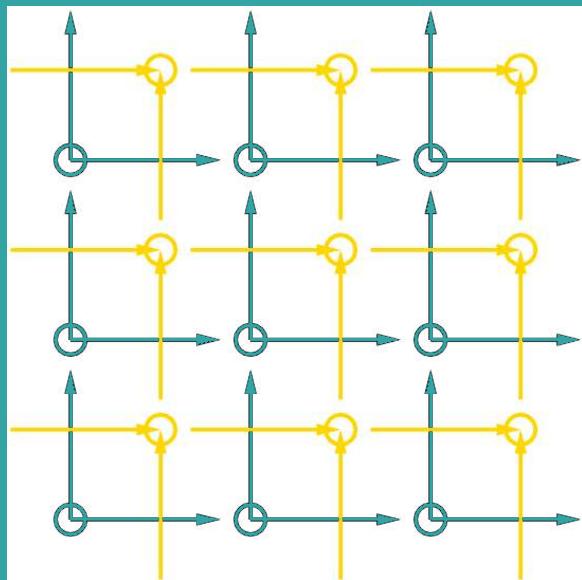
# What is computed



- Resonant fields
  - Periodic boundary conditions
- Time dependent fields excited via
  - Relativistic line charges
  - Free moving charges
  - Portmodes

Materials can be lossy and dispersive

# Finite Difference/ Finite Integration



- Electric/ magnetic voltages at the edges of cells

$$\begin{aligned}-\frac{d}{dt}h &= \int_{\Delta s'} \frac{1}{\int_A \mu dA} ds' (e_1 + e_2 - e_3 - e_4) \\ \frac{d}{dt}e &= \int_{\Delta s} \frac{1}{\int_A \varepsilon dA'} ds (h_1 + h_2 - h_3 - h_4)\end{aligned}$$

# FD-Equations

Loss free Maxwell

$$\int \frac{d}{dt} \varepsilon \vec{E} \cdot d\vec{A} = \oint \vec{H} \cdot d\vec{s}$$

$$- \int \frac{d}{dt} \mu \vec{H} \cdot d\vec{A} = \oint \vec{E} \cdot d\vec{s}$$

We approximate

$$\int \varepsilon dA' \frac{d}{dt} \vec{E} \cdot \vec{n}'_A \approx \oint \vec{H} \cdot d\vec{s}'$$

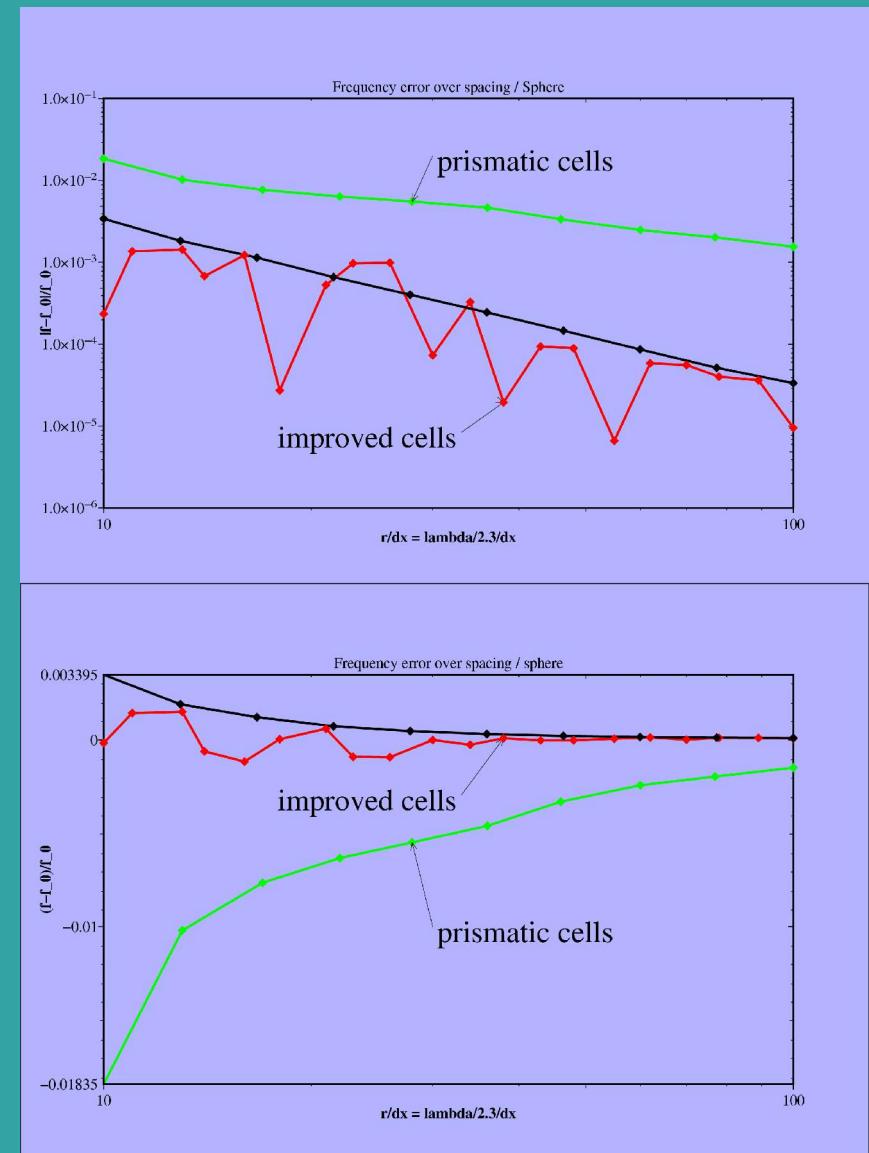
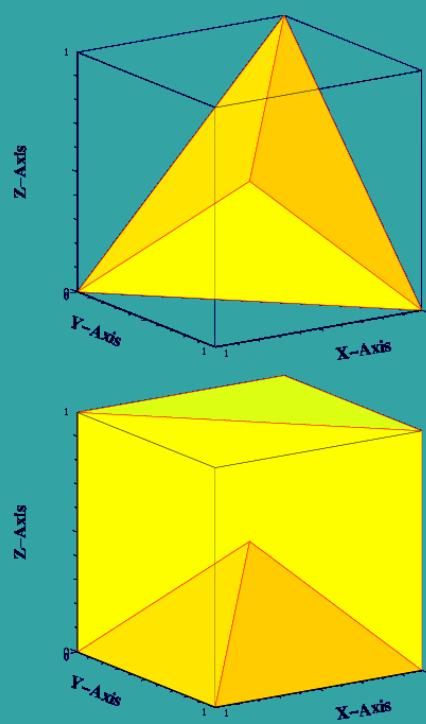
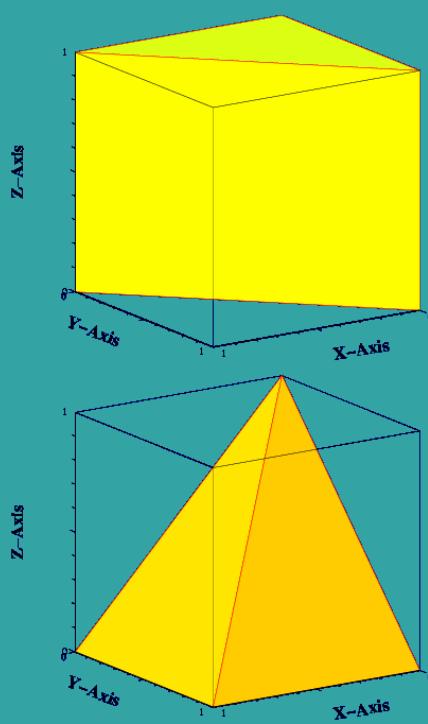
$$- \int \mu dA \frac{d}{dt} \vec{H} \cdot \vec{n}_A \approx \oint \vec{E} \cdot d\vec{s}$$

Integration gives electric/  
magnetic voltages

$$\frac{d}{dt} \int \vec{E} \cdot d\vec{s} \approx \frac{\int ds}{\int \varepsilon dA'} \oint \vec{H} \cdot d\vec{s}'$$

$$- \frac{d}{dt} \int \vec{H} \cdot d\vec{s}' \approx \frac{\int ds'}{\int \mu dA} \oint \vec{E} \cdot d\vec{s}$$

# Generalised diagonal fillings

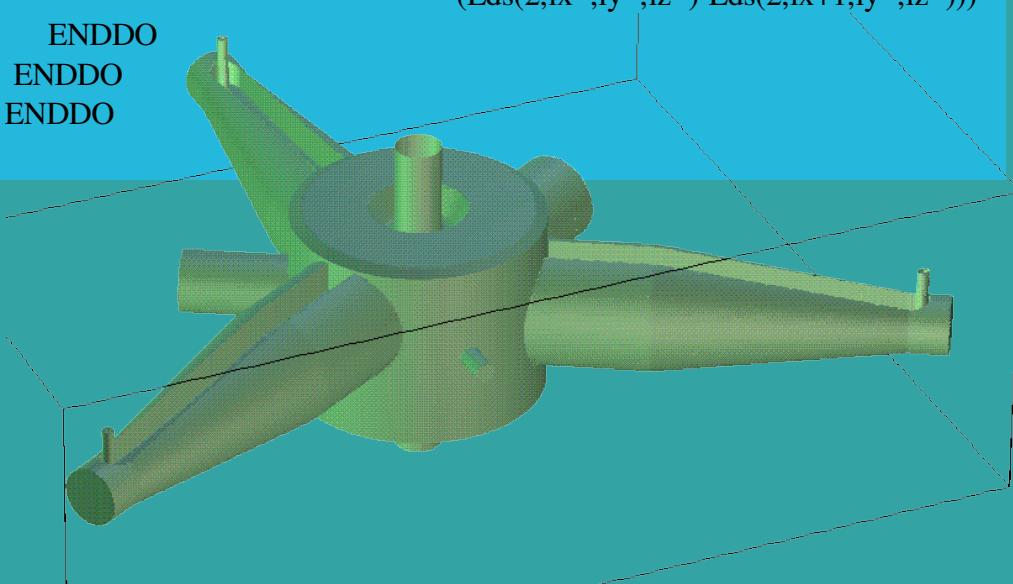


# Memory and CPU-Time only for interesting Gridcells

```

REAL, DIMENSION(1:3,0:nx+1,0:ny+1,0:nz+1) :: &
    Eds, Hds, dsoEpsA, dsoMuA
DO iz= 1, nz, 1
    DO iy= 1, ny, 1
        DO ix= 1, nx, 1
            Hds(1,ix,iy,iz)= Hds(1,ix,iy,iz) &
                - dt*dsoMuA(1,ix,iy,iz) * ( Eds(2,ix ,iy ,iz )-Eds(2,ix ,iy ,iz+1) &
                    + Eds(3,ix ,iy ,iz )-Eds(3,ix ,iy+1,iz ))
            Hds(2,ix,iy,iz)= Hds(2,ix,iy,iz) &
                - dt*dsoMuA(2,ix,iy,iz) * (-Eds(1,ix ,iy ,iz )-Eds(1,ix ,iy ,iz+1)) &
                    + Eds(3,ix ,iy ,iz )-Eds(3,ix+1,iy ,iz ))
            Hds(3,ix,iy,iz)= Hds(3,ix,iy,iz) &
                - dt*dsoMuA(3,ix,iy,iz) * ( Eds(1,ix ,iy ,iz )-Eds(1,ix ,iy+1,iz ) &
                    -(Eds(2,ix ,iy ,iz )-Eds(2,ix+1,iy ,iz )))
        ENDDO
    ENDDO
    ENDDO

```

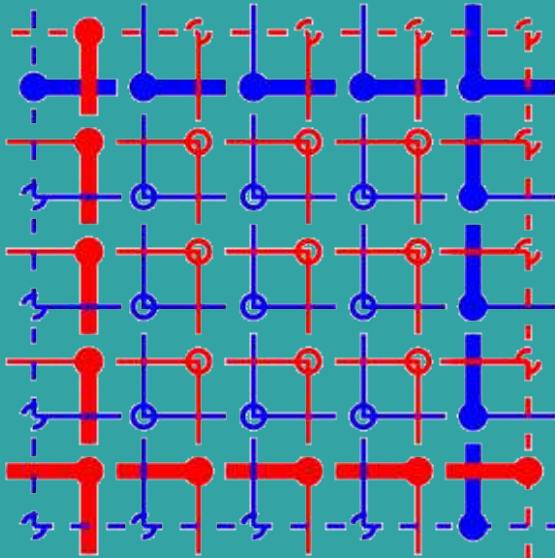


```

REAL,DIMENSION(1:3,0:NetCells) :: Eds, Hds
INTEGER, DIMENSION(0:NNetCells) :: KindofCell
INTEGER, DIMENSION(0:nx+1,0:ny+1,0:nz+1) :: NrofCell
REAL, DIMENSION(1:3,0:NDifferentCells) :: dsoEpsA, dsoMuA
DO iz= 1, nz, 1
    DO iy= 1, ny, 1
        DO ix= 1, nx, 1
            i= NrofCell(ix,iy,iz)
            IF ( i .LT. 1) CYCLE
            k= KindofCell(i)
            ipx= NrofCell(ix+1,iy ,iz )
            ipy= NrofCell(ix ,iy+1,iz )
            ipz= NrofCell(ix ,iy ,iz+1)
            Hds(1,i)= Hds(1,i) - dt*dsoMuA(1,k) &
                * ( Eds(2,i)-Eds(2,ipz) + Eds(3,i)-Eds(3,ipy) )
            Hds(2,i)= Hds(2,i) - dt*dsoMuA(2,k) &
                * (-Eds(1,i)-Eds(1,ipz)) + Eds(3,i)-Eds(3,ipx) )
            Hds(3,i)= Hds(3,i) - dt*dsoMuA(3,k) &
                * ( Eds(1,i)-Eds(1,ipy) -(Eds(2,i)-Eds(2,ipx)))
        ENDDO
    ENDDO
    ENDDO

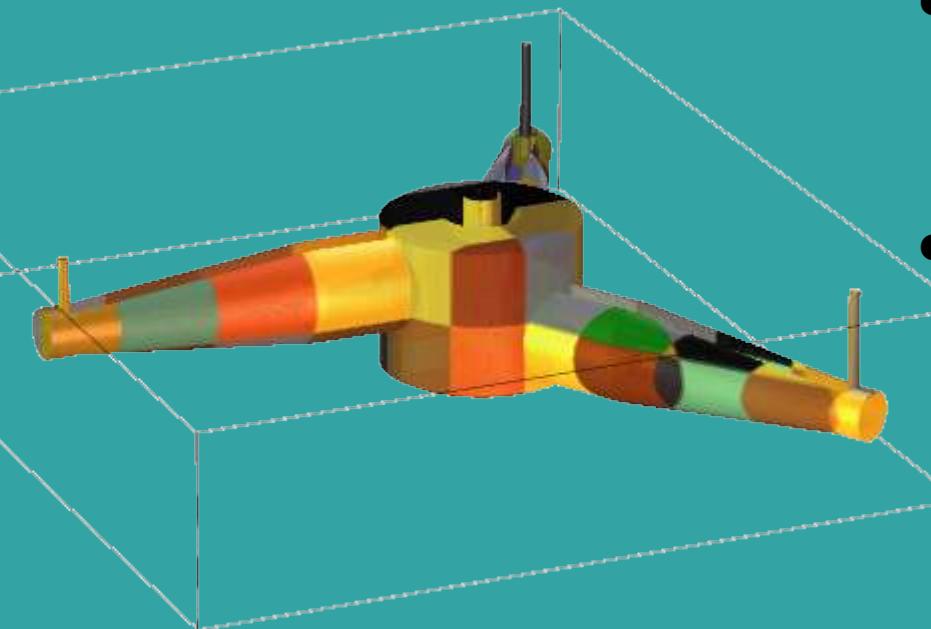
```

# Parallel Computation



For all Timesteps: DO  
    Compute local H by applying the local curl operator to the local E  
        For all Directions: DO  
            Send tangential H to the neighbour  
            Receive tangential H from the neighbour  
        ENDDO For all Directions  
    Compute local E  
        For all Directions: DO  
            Send tangential E to the neighbour  
            Receive tangential E from the neighbour  
        ENDDO For all Directions  
ENDDO For all Timesteps

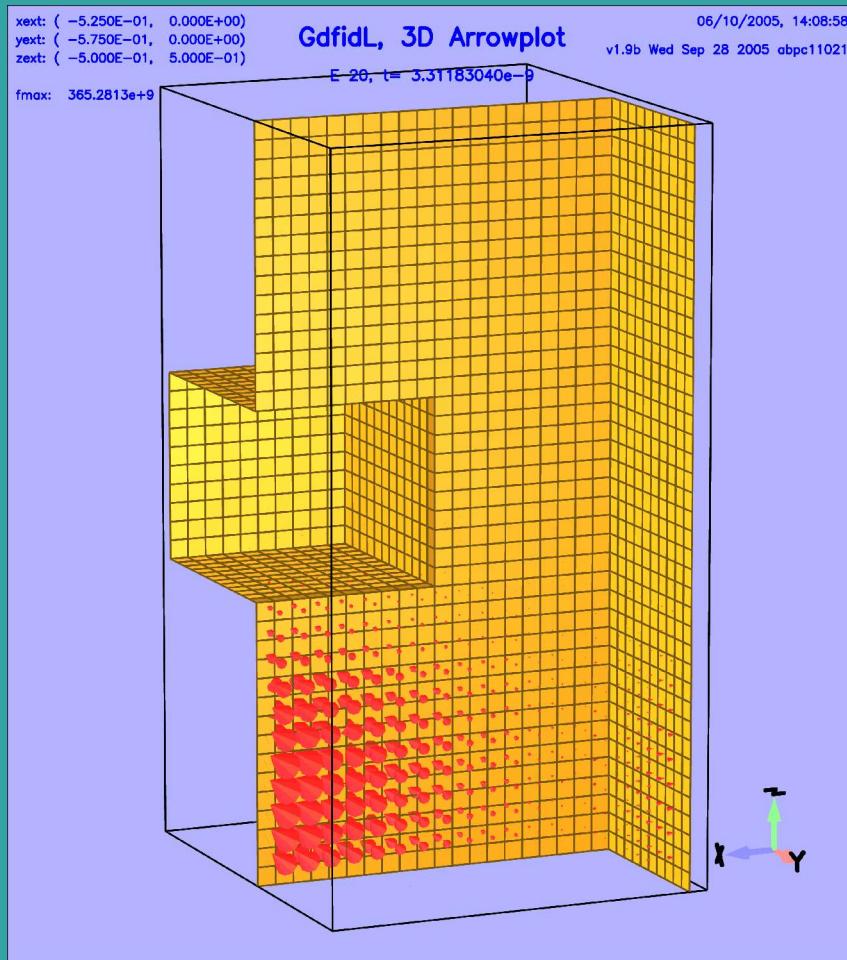
# Parallel Computation, Load Balancing



- Volume is partitioned in many subvolumes
- Interesting subvolumes are spread over the available processors.

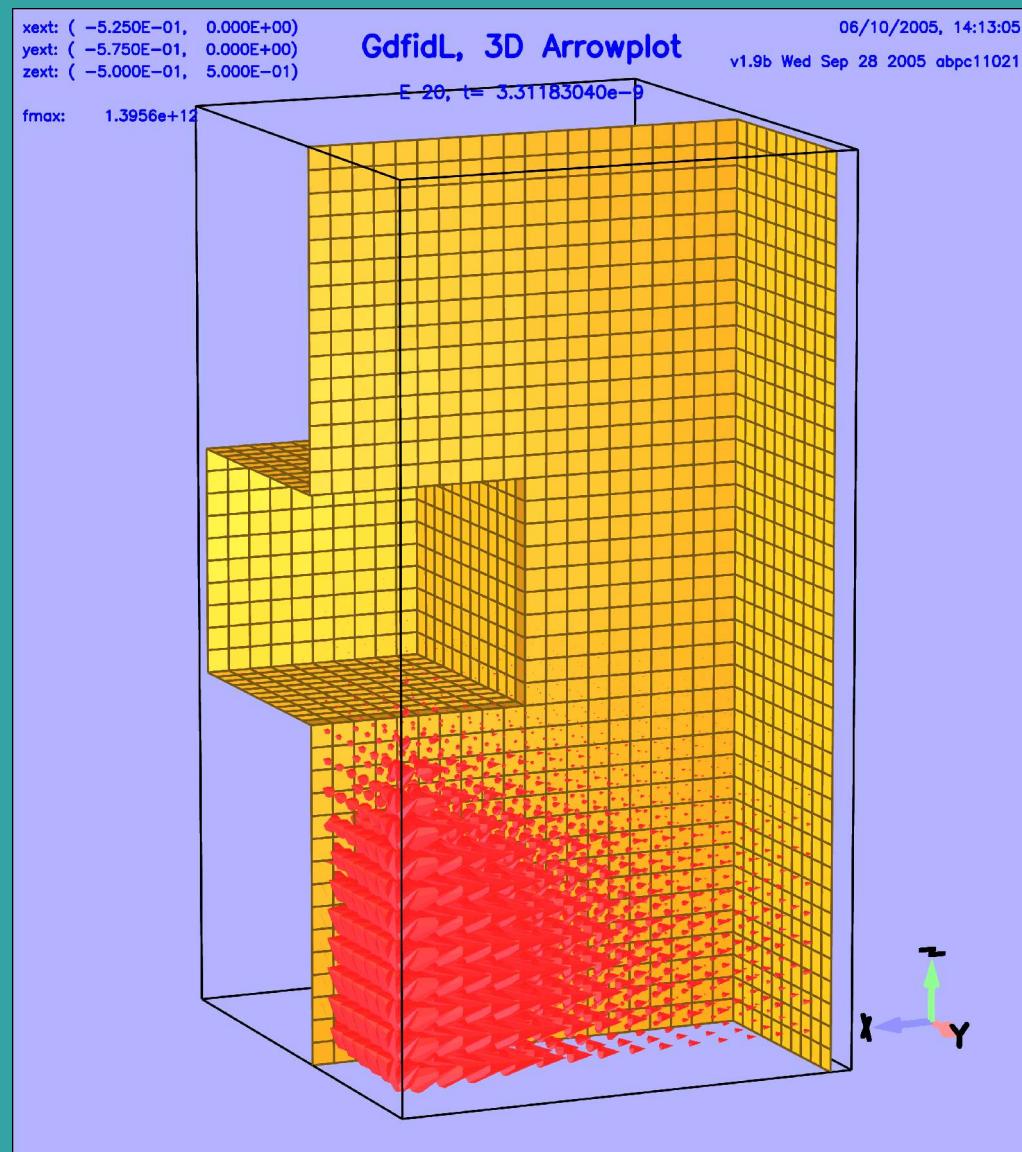
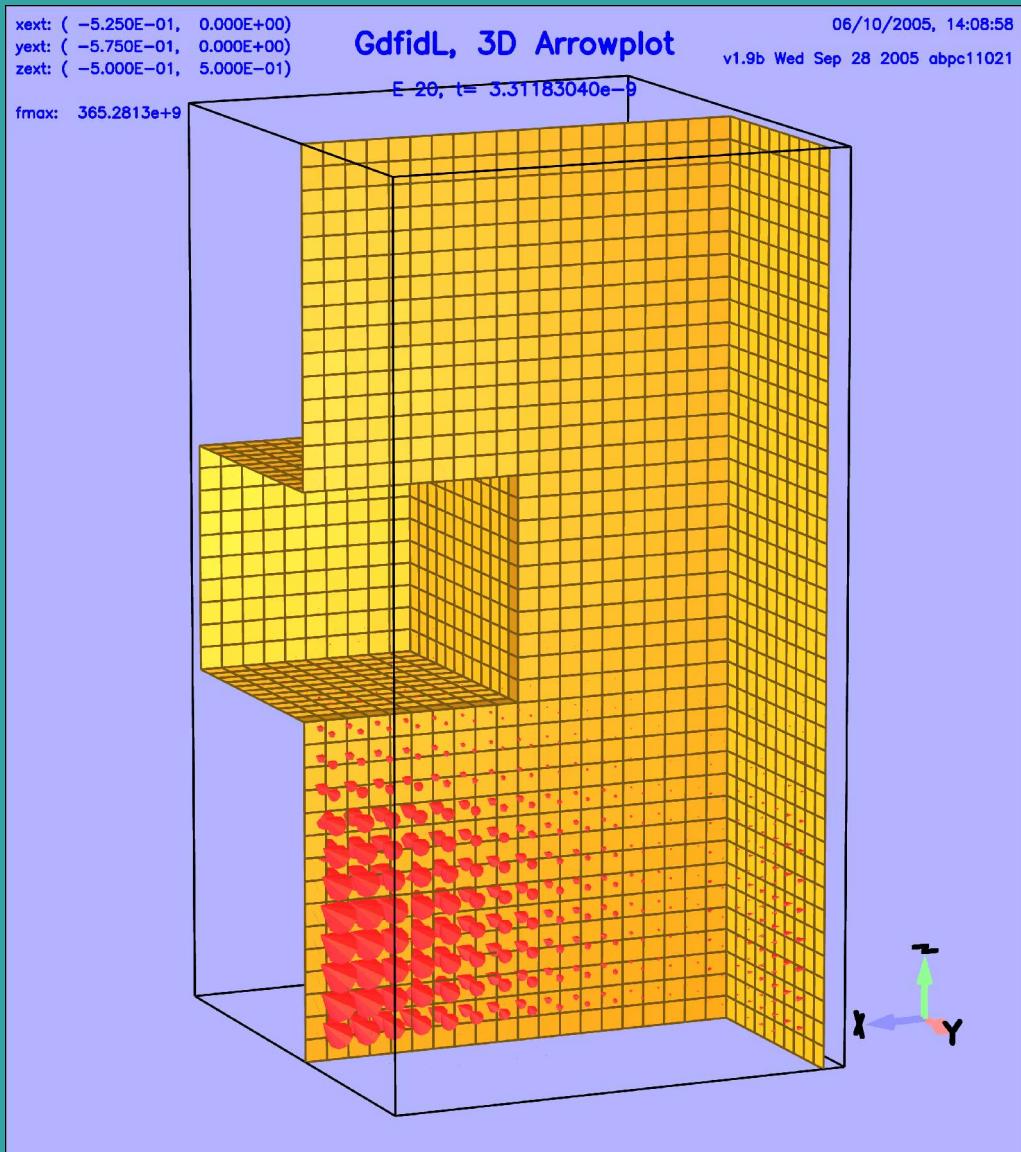
Less than 10% interesting volume

# Excitation with hollow Charge

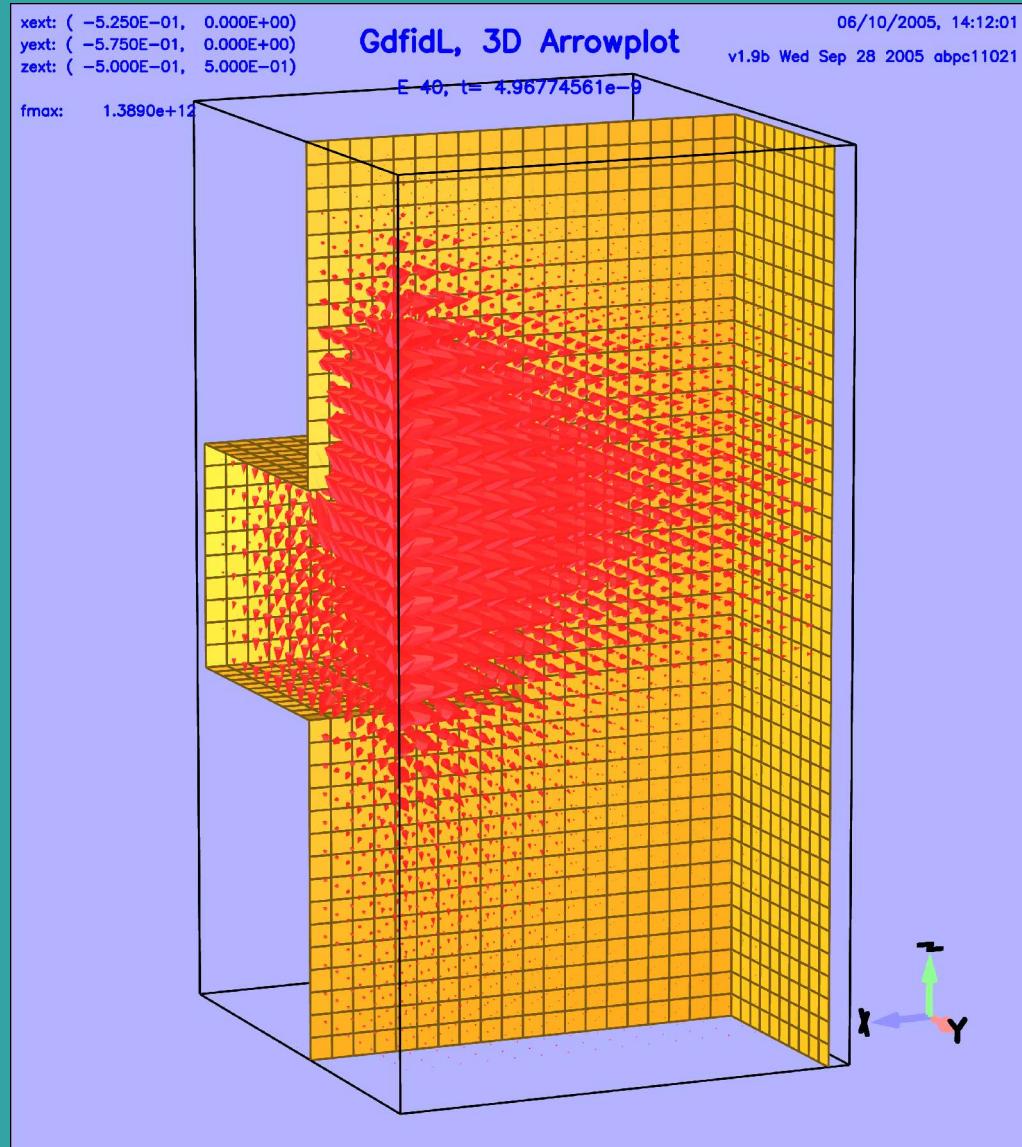
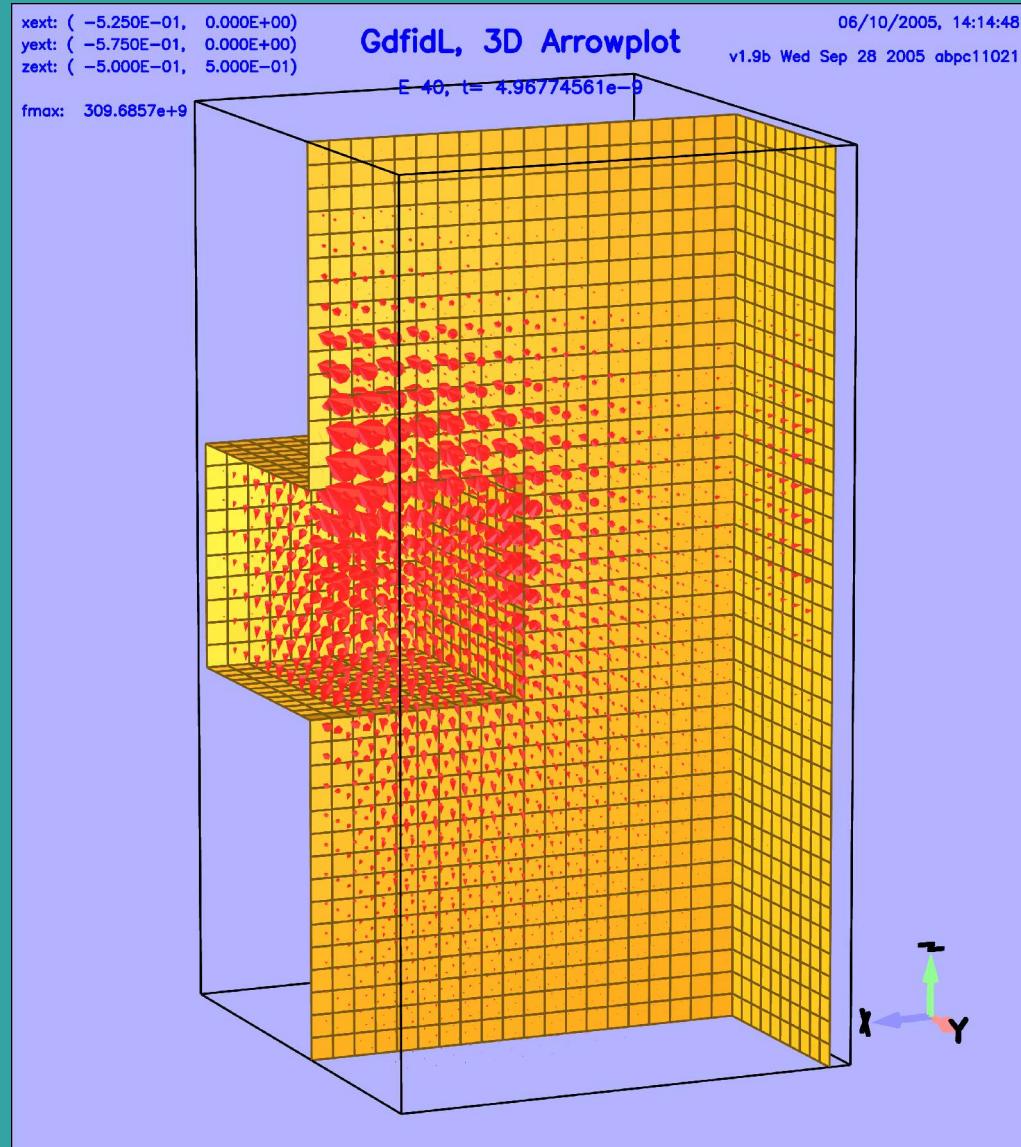


- Hollow charge very near to the beampipe
- Less dispersion error

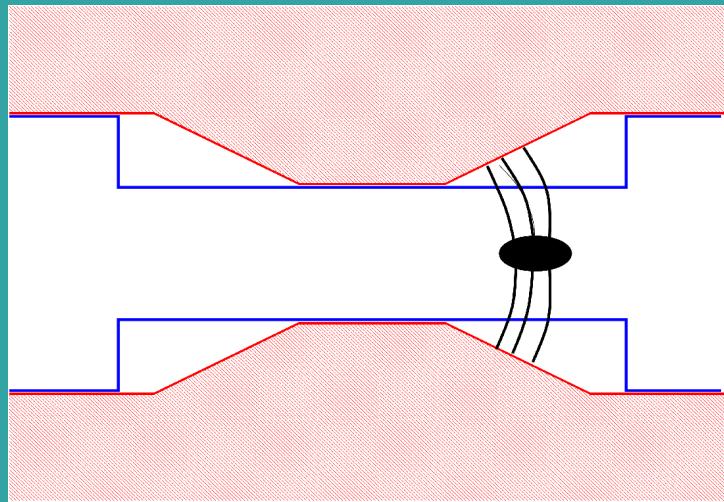
# Hollow Charge



# Hollow Charge



# Napoly Integration in 3D



- Avoids "Catch-up Problem" also for Collimator-like structures.

$$W_z(x, y, s) = \frac{c}{Q} \int_{-\infty}^{\infty} E_z(x, y, z = ct - s, t) dt$$

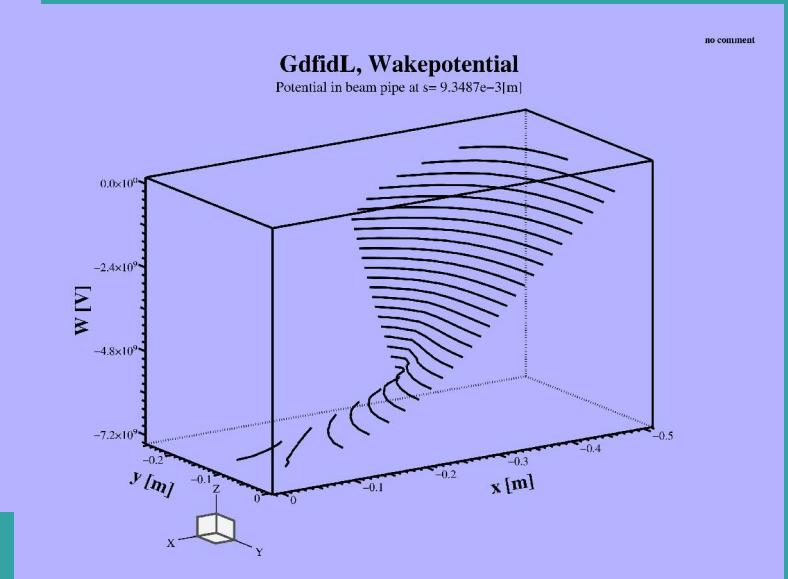
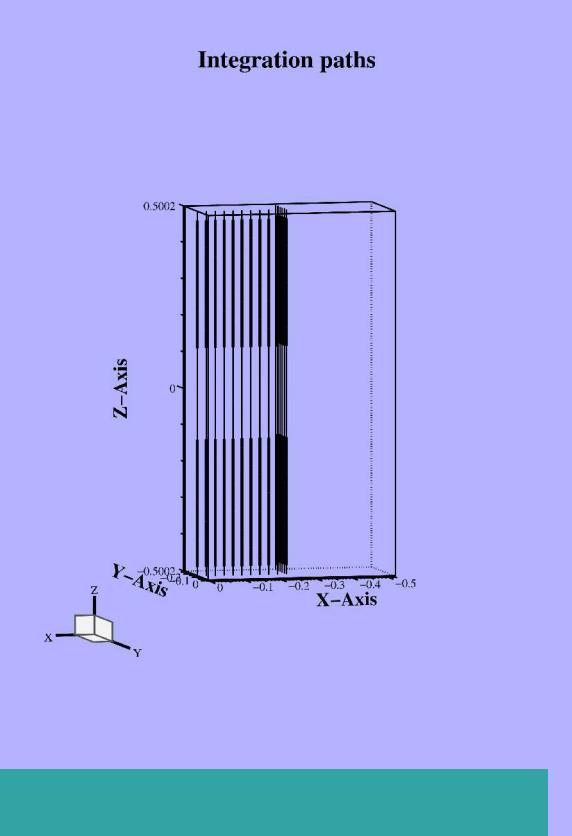
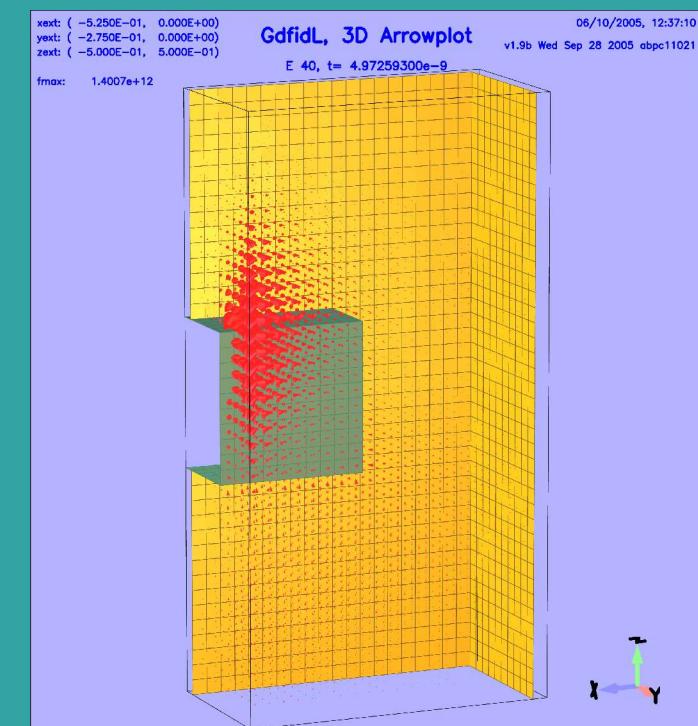
$$c \int_{t=s/c}^{\infty} E_z(x, y, z = ct - s, t) dt = \int_{\xi=x}^{x_{Max}} (E_x^{TM} + cB_y^{TM}) (x = \xi, y, z = 0, t = s/c) d\xi$$

# Napoly Integration in 3D

- Integration only over TM part of the field.
- Requires solution of 2D Poisson problems at each timestep

$$\begin{aligned}\vec{E}_t^{TM} &= \nabla_t(\varphi^{TM}) \\ \Delta_t \varphi^{TM} &= -\frac{d}{dz} E_z \\ \vec{H}_t^{TM} &= \vec{H}_t - \vec{H}_t^{TE} \\ \vec{H}_t^{TE} &= \nabla_t(\varphi^{TE}) \\ \Delta \varphi^{TE} &= -\frac{d}{dz} H_z\end{aligned}$$

# Napoly Integration



# Summary

- Computation on parallel systems
  - Incredibly large number of gridcells possible
- Wakepotentials
  - Napolyn Integration for general geometries
  - Hollow beam excitation for general geometries
- Generalised diagonal fillings
- Eigenvalues of lossy structures