

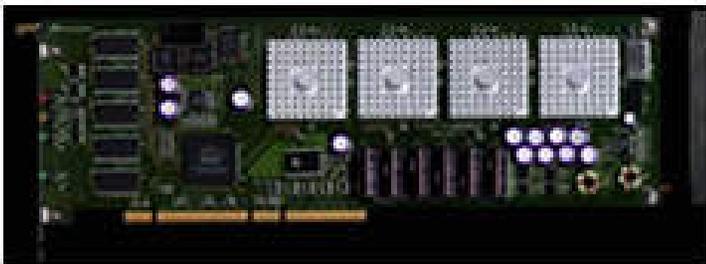
Development of a Fast Vortex Method for Fluid Flow Simulation using Special- Purpose Computers

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1

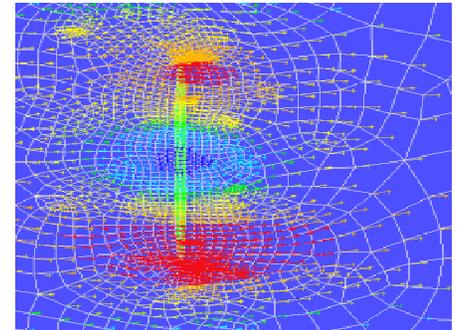
Introduction

1.1 Background

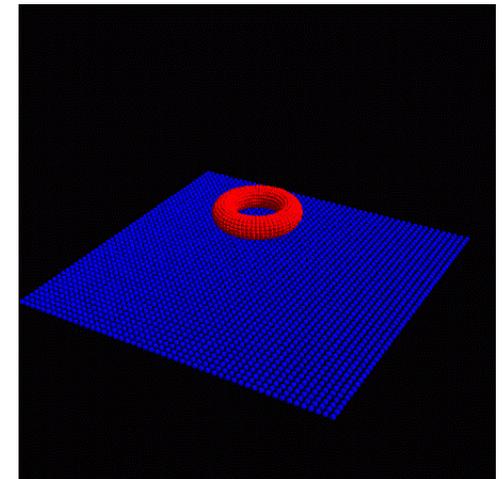
- Vortex methods introduced in 1930s by Rosenhead
- Developed for complicated, unsteady and vortical flows
- N-Body algorithm introduced in 1950s
- Digital computers introduced in 1970s
- N-Body simulation is particle-interactions and calculation cost is $O(N^2)$
- Vortex method is one of the N-body algorithms

1.1 Background(Contd...)

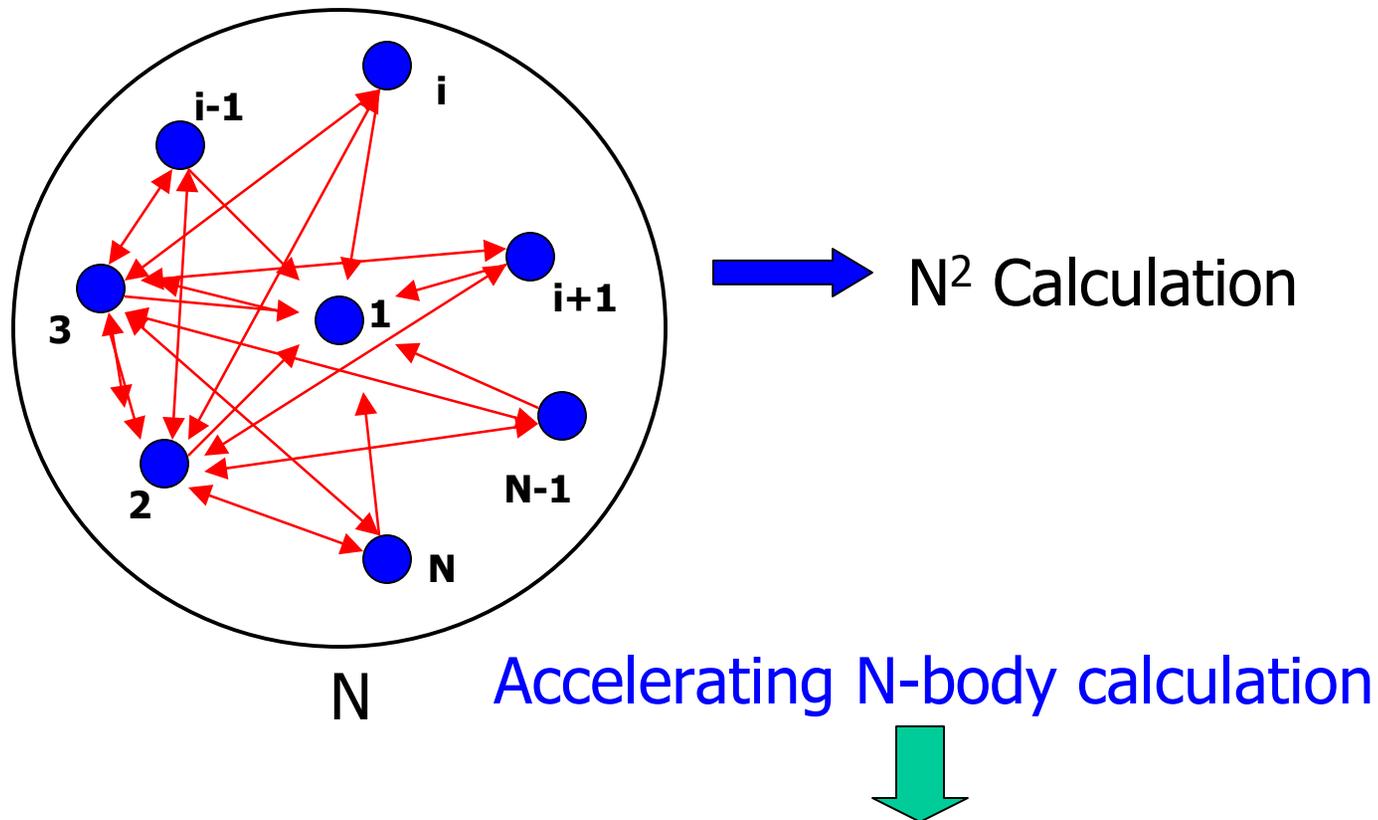
- Advantages of Vortex Method (VM)
 - Lagrangian based CFD method
 - Calculates only regions of non-zero vorticity
 - Can be applied to high Reynolds number flows in complex geometry
 - Can be solved convection in straightforward
- Disadvantages of VM
 - High computation cost
 - Descretization error
 - Diffusion error



Asakura, 2002



1.2 The need for acceleration techniques



Hardware: parallel calculation

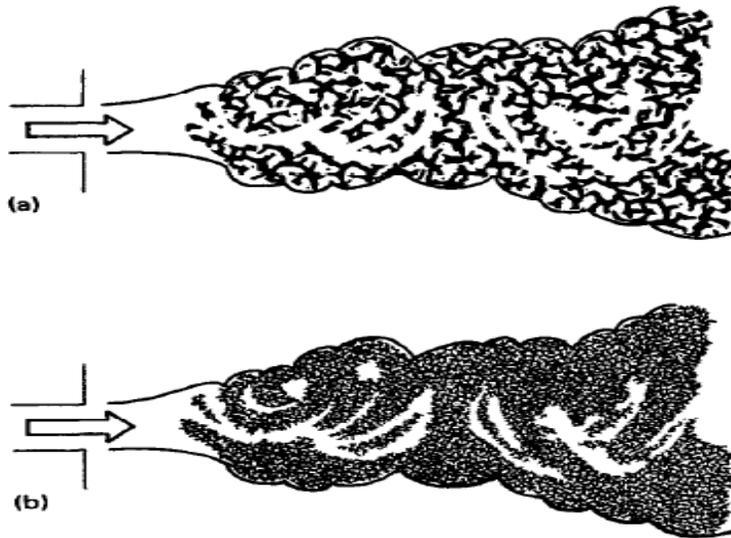
Software: Tree code, FMM, P^2M^2 Pseudo-particle multipole method etc.

1.3 Motivation

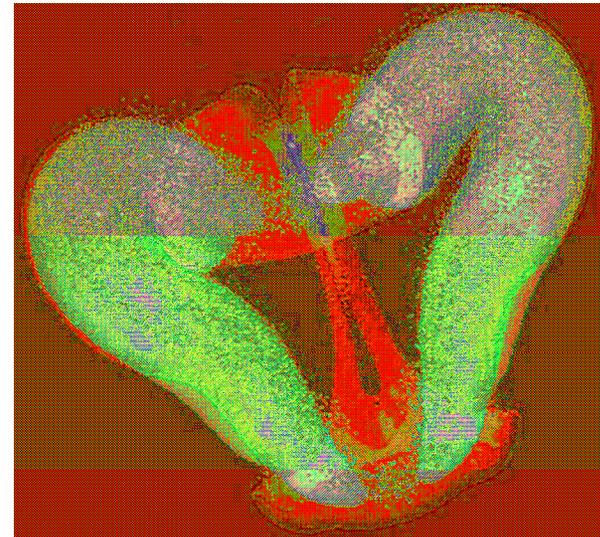
- Two ways to reduce calculation cost
 - Fast algorithms (Cheng, 1999)
 - Special-purpose computers (Susukita, 2003; Narumi, 2006)
- Fast algorithm has high proportionality cost at higher accuracy (Greengard, 1987)
- Special-purpose computers have been developed to accelerate MD simulations (Narumi, 1997)
 - It can be applied to accelerate VM calculation
 - Fast algorithms can be implemented
 - It does not support for Fast Poisson solver
- Recently GPGPU has been used to accelerate VM calculation (Stock, 2008)

1.3 Motivation(Contd...)

- To calculate for high Reynolds number turbulent flows which required high performance computational resources



Tennekes and Lumely, A first course
In turbulence



Kida, 1994

- The collision of vortex rings contain millions of particles result in a highly turbulent state

1.4 Previous Studies

- Vortex rings have been studied in the broader arena of vortex interaction (Shariff, 1992)
- Large N is necessary to capture the essential characteristics of vortex rings collisions (Winckelmans, 1993)
- High Reynolds number is necessary to generate a secondary vortex rings (Mammetti, 1999)
- Computational resources are essential for longer calculation and to produce the fast mechanism of energy transfer (Chatelain, 2003)
- Fast Poisson solvers are still faster compared with VIC method (Cottet, 2002)
- Fast algorithm is successfully implemented on special-purpose computers for astrophysical problems (Makino, 1991, Kawai, 2004)

1.5 Purpose of the present study

- To accelerate the high Reynolds number VM calculation without loss of numerical accuracy
- To develop a fast vortex method using special-purpose computers
- To solve the three critical issues
 - The efficient calculation of Biot-Savart law and stretching term
 - An optimized function table
 - Round-off error caused by the partially single precision of MDGRAPE
 - Special treatments for cross product calculation
- To implement fast algorithms on special-purpose computers for further acceleration
- Comparative study to validate this scheme

2

Numerical Methods

2.2 Vortex Methods

- Lagrangian methods used to simulate unsteady, convection-dominated problems
- It has a difficulty in achieving higher order spatial accuracy compared to Eulerian methods
- It is required to consider an accurate viscous diffusion scheme

2.2.1 Formulation of 3D VM

- Vorticity transport:
$$\frac{D\boldsymbol{\omega}}{Dt} = (\boldsymbol{\omega} \cdot \nabla)\mathbf{u} + \nu \nabla^2 \boldsymbol{\omega} \quad (2.1)$$

- Biot-Savart Law:
$$\mathbf{u}(\mathbf{x}) = -\frac{1}{4\pi} \int \frac{(\mathbf{x} - \mathbf{x}') \times \boldsymbol{\omega}(\mathbf{x}')}{|\mathbf{x} - \mathbf{x}'|^3} dV(\mathbf{x}') \quad (2.2)$$

- Discretized form:
$$\mathbf{u}_i = -\frac{1}{4\pi} \sum_{j=1}^N \frac{\mathbf{r}_{ij}^2 + (5/2)\sigma_j^2}{(\mathbf{r}_{ij}^2 + \sigma_j^2)^{5/2}} \mathbf{r}_{ij} \times \boldsymbol{\gamma}_j \quad (2.3)$$

- Stretching term:
$$\frac{d\boldsymbol{\gamma}_i}{dt} = \frac{1}{4\pi} \sum_{j=1}^N \left\{ -\frac{\mathbf{r}_{ij}^2 + (5/2)\sigma_j^2}{(\mathbf{r}_{ij}^2 + \sigma_j^2)^{5/2}} \boldsymbol{\gamma}_i \times \boldsymbol{\gamma}_j + 3 \frac{\mathbf{r}_{ij}^2 + (7/2)\sigma_j^2}{(\mathbf{r}_{ij}^2 + \sigma_j^2)^{7/2}} (\boldsymbol{\gamma}_i \cdot \mathbf{r}_{ij})(\mathbf{r}_{ij} \times \boldsymbol{\gamma}_j) \right\} \quad (2.6)$$

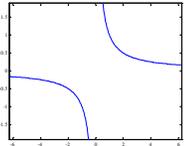
Core Spreading Method

- Viscous Diffusion:

$$\frac{d\omega_i}{dt} = \nu \nabla^2 \omega_i \quad (2.7)$$

- Green's Function Solution:

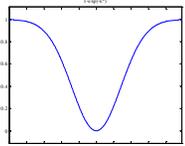
$$\omega_i = \frac{\gamma_j}{(4\pi\nu t)^{d/2}} \exp\left(-\frac{|\mathbf{x}_j - \mathbf{x}_i|^2}{4\nu t}\right) \quad (2.8)$$



- Vorticity at arbitrary point:

$$\omega_i = \sum_j \gamma_j \zeta(|\mathbf{x}_j - \mathbf{x}_i|) \quad (2.9)$$

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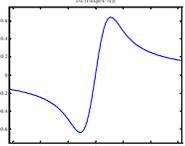
- Cutoff Function:

$$\zeta = \frac{1}{(2\pi\sigma^2)^{d/2}} \exp\left(-\frac{|\mathbf{x}_j - \mathbf{x}_i|^2}{2\sigma^2}\right) \quad (2.10)$$

||

- Core expansion rate:

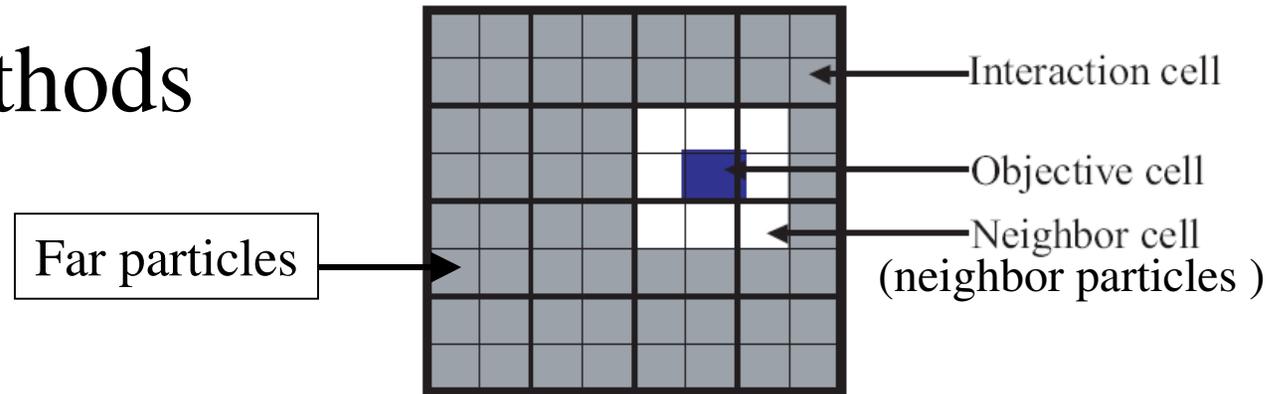
$$\sigma^2 = 2\nu t \quad (2.11)$$



- Position update:

$$\frac{d\mathbf{x}_i}{dt} = \mathbf{u}_i \quad (2.12)$$

2.3 Fast Methods



- Tree Algorithm (Barnes and Hut, 1986)
 - Hierarchical data structure
 - Calculation cost $O(N \log N)$
 - Can be implemented on special-purpose computers
- Fast Multipole Method (Greengard and Rokhlin, 1987)
 - All particles are uniformly distributed in a unit cube
 - Far particles calculated as a multipole expansion
 - Neighbor particles calculated in a direct summation
 - Calculation cost is proportional to $O(N)$
- Other fast methods
 - Anderson's method (Anderson, 1992)
 - Pseudo-particle multipole method (Makino, 1999)

2.3 Fast Methods(Contd...)

- FMM has been used in my calculation
- Biot-Savart equation has been derived as

$$\mathbf{u}_i \approx \frac{1}{4\pi} \sum_{n=0}^p \sum_{m=-n}^n \left\{ \sum_{j=1}^N \gamma_j M_j \right\} \times \nabla S_i \quad (2.13)$$

$$\mathbf{u}_i \approx \frac{1}{4\pi} \sum_{n=0}^p \sum_{m=-n}^n \left\{ \sum_{j=1}^N \gamma_j L_j \right\} \times \nabla R_i \quad (2.14)$$

- Stretching term derived as

$$\frac{D\gamma_i}{Dt} \approx \frac{1}{4\pi} \sum_{n=0}^p \sum_{m=-n}^n \left\{ \sum_{j=1}^N \gamma_j \times \nabla M_j \right\} (\gamma_i \cdot \nabla S_i) \quad (2.15)$$

$$\frac{D\gamma_i}{Dt} \approx \frac{1}{4\pi} \sum_{n=0}^p \sum_{m=-n}^n \left\{ \sum_{j=1}^N \gamma_j \times \nabla L_j \right\} (\gamma_i \cdot \nabla R_i) \quad (2.16)$$

Here p is order of moment and L, M, R, S are denoted to simplify the equations.

Summary

- A fast vortex method has been formulated
- 3D core spreading method is explained
- Different fast methods are reviewed briefly

4

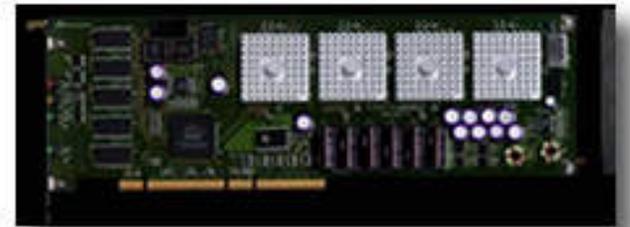
Fast Vortex Method Calculation
using a Special-purpose
Computer

4.1 Introduction

- A mathematical formulation will be developed for VM calculation using MDGRAPE-2
- The efficient calculation of Biot-Savart law and stretching term will be performed
- An optimum range of a function table is determined
- The cross-product calculation will be demonstrated
- The accuracy will be evaluated by calculating the impingement of two identical vortex rings
- The calculated results is compared with and without the use of MDGRAPE-2 and with referenced work carefully

Special-purpose Computer: MDGRAPE-2 (Narumi, 1997)

- PCI-board for accelerating Molecular Dynamics Simulation
- Reduced computation cost significantly for N^2 calculation
- Particle Memory: 5.0×10^5 (20MB)
- Calculation speed: 64 Gflops
- Speed up 10-1000 times faster
- Compatible for FORTRAN and C programming Languages



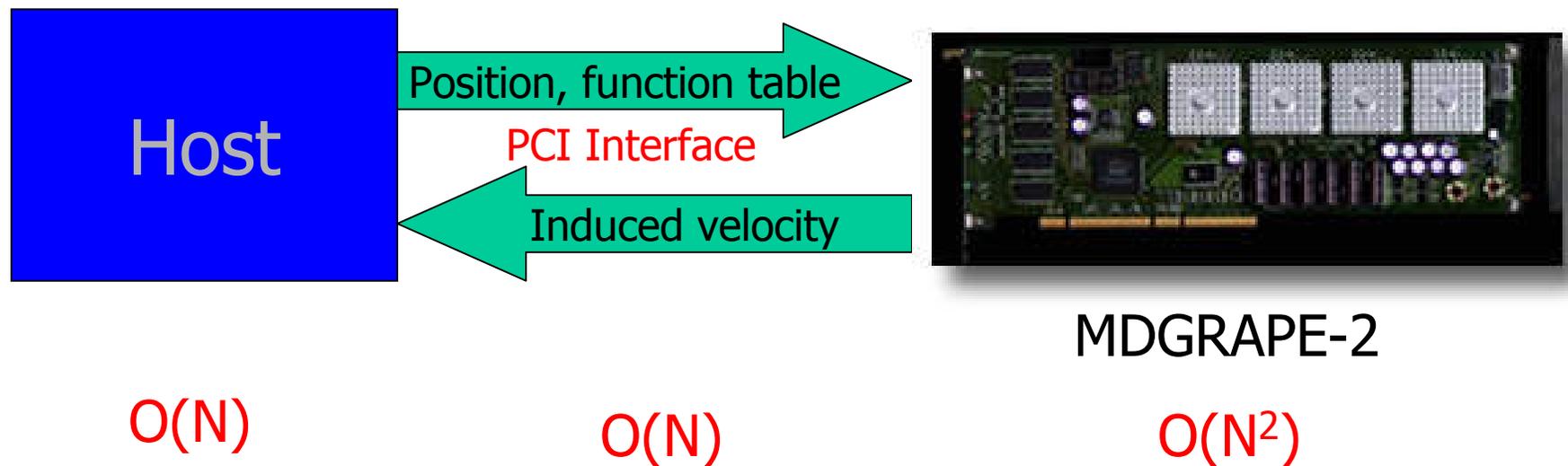
Host Machine and MDGRAPE-2

General-purpose

Time Integration, Vorticity, etc..

Special-purpose

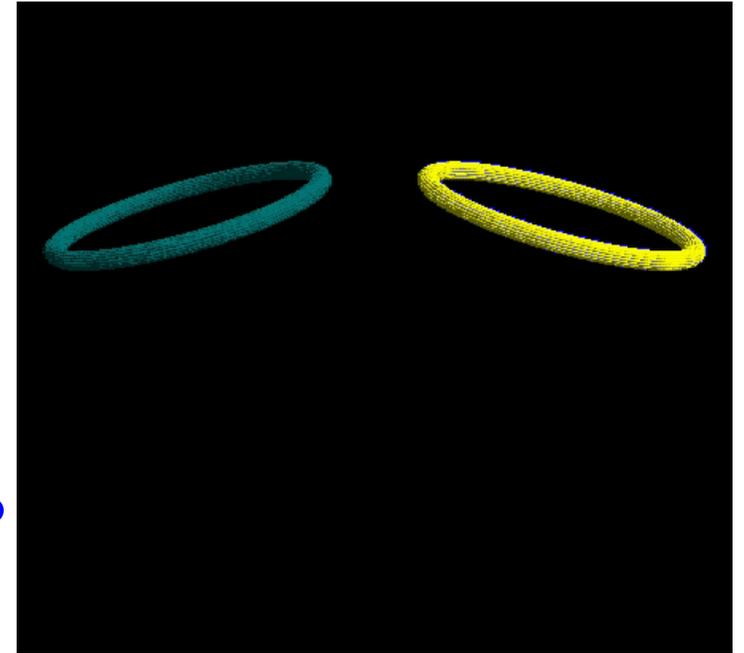
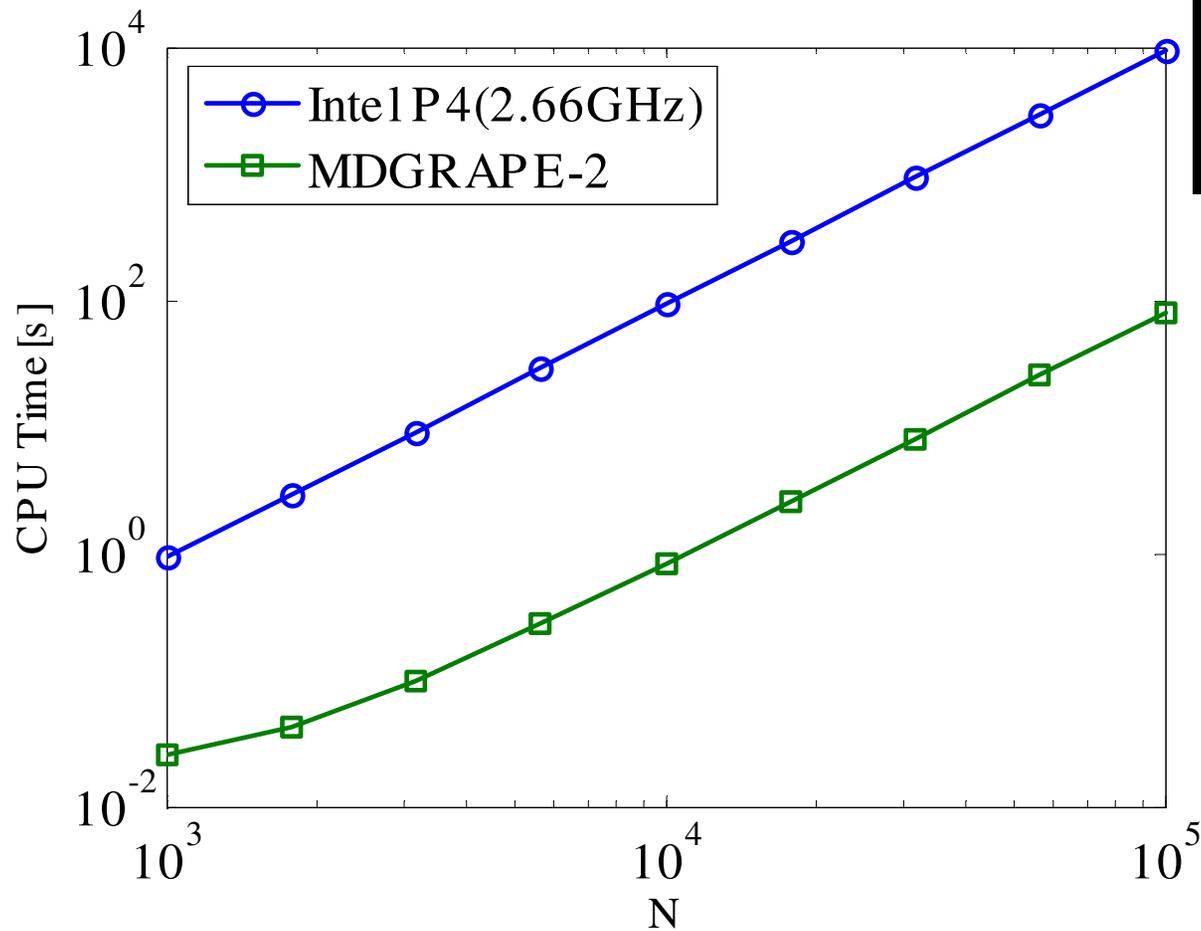
Induced Velocity, Stretching Term



Hosts for performance measurement

Special-purpose	Host CPU	Cache	Memory	OS	Compiler
MDGRAPE-2	Intel P4 2.66GHz (1CPU 1 Core)	512KB	1GB (2 GB Swap Memory)	Linux 8.0 Kernel 2.4.18-14	ifort
MDGRAPE-3	Xeon 5160 3.0GHz (1CPU 2 Core)	4096KB	32GB (0 GB Swap Memory)	Cent 4.3 (Final) Kernel 2.6.9- 34_ELsmp	ifort

Performance



$N=6 \times 10^4$

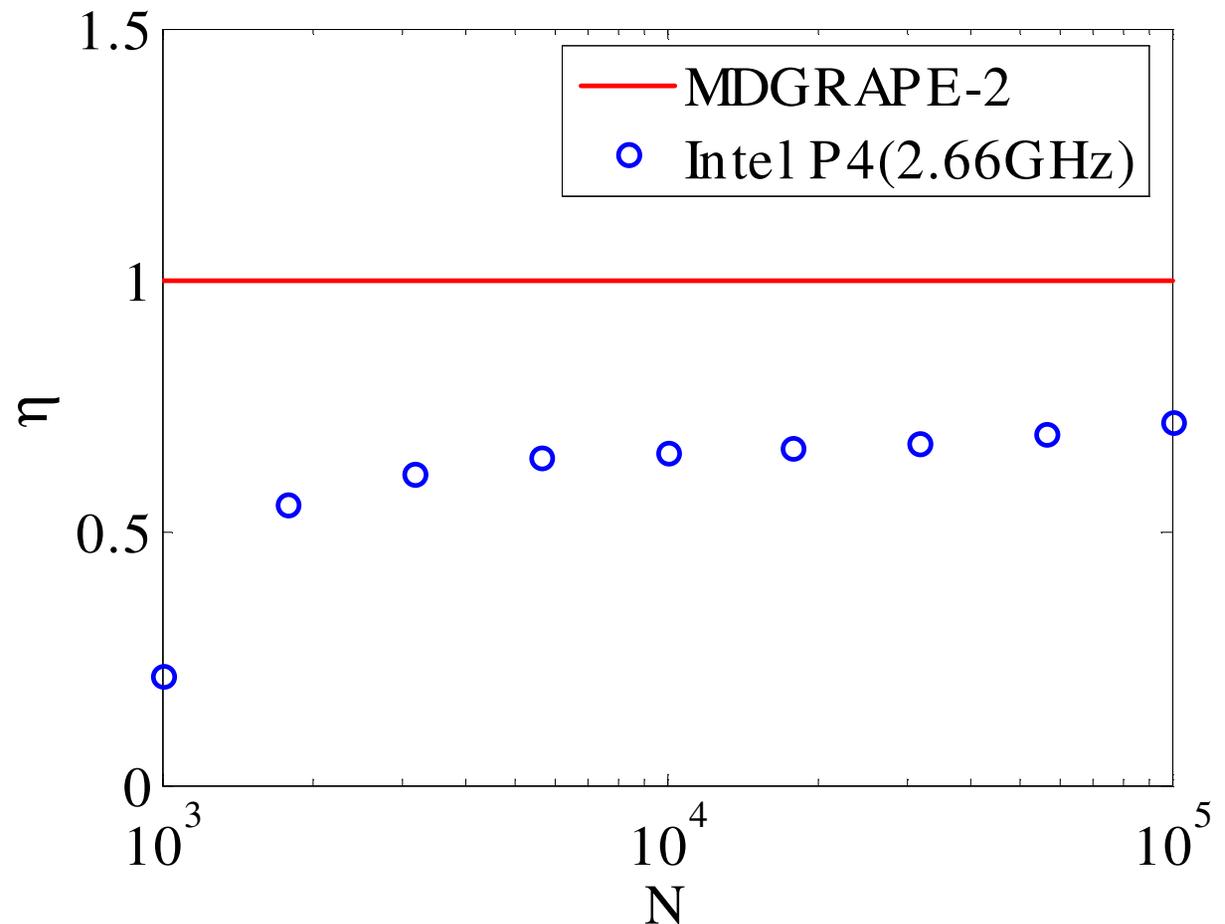
100 times at $N \sim 10^5$

Efficiency of VM calculation

$$N_{APPL} = \frac{nmd(xmd + ymd + zmd)}{CPUtime(sec/step)}; \quad N_{GRAPE} = 5.5 \times 10^8$$

Efficiency:

$$\eta = \frac{N_{APPL}}{N_{GRAPE}}$$



Difficulties with MDGRAPE-2 for VM

- Not designed for vector product calculation
- Requires optimum generation of a function table
- Calculates partly with single precision

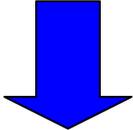
Special treatments are required to solve these problems

4.2 Mathematical Formulations

Coulomb Potential:
$$\Phi_i = \sum_{j=1}^N b_{ij} g(w) \mathbf{r}_{ij} = \sum_{j=1}^N b_{ij} g\left(a_{ij} \left(|\mathbf{r}_{ij}|^2 + \epsilon_{ij}^2\right)\right) \quad (4.1)$$

Coulomb Force:
$$\mathbf{f}_i = \sum_{j=1}^N b_{ij} g(w) \mathbf{r}_{ij} = \sum_{j=1}^N b_{ij} g\left(a_{ij} \left(|\mathbf{r}_{ij}|^2 + \epsilon_{ij}^2\right)\right) \mathbf{r}_{ij} \quad (4.2)$$

Vortex Method (Induced velocity)

$$\mathbf{u}_i = \sum_{j=1}^N \frac{\mathbf{r}_{ij}^2 + (5/2)\sigma_j^2}{4\pi \left(\mathbf{r}_{ij}^2 + \sigma_j^2\right)^{5/2}} \mathbf{r}_{ij} \times \gamma_j$$

$$\mathbf{u}_i = \sum_{j=1}^N B_j g\left(A_j \left(|\mathbf{r}_{ij}|^2 + \epsilon_{ij}^2\right)\right) \mathbf{r}_{ij} \quad (4.3)$$

4.2 Mathematical Formulations (contd.)

- Input

$$\mathbf{r}_{ij} = (x_{ij}, y_{ij}, z_{ij}) \quad \boldsymbol{\gamma}_j = (\gamma_j^x, \gamma_j^y, \gamma_j^z) \quad (4.4)$$

- Vector product

$$\sum_j \mathbf{r}_{ij} \times \boldsymbol{\gamma}_j = \sum_j (y_{ij} \gamma_j^z - z_{ij} \gamma_j^y, z_{ij} \gamma_j^x - x_{ij} \gamma_j^z, x_{ij} \gamma_j^y - y_{ij} \gamma_j^x) \quad (4.5)$$

- 3 x Scalar product

$$\sum_j \mathbf{r}_{ij} \cdot \boldsymbol{\gamma}_j^x = \sum_j (x_{ij} \gamma_j^x, y_{ij} \gamma_j^x, z_{ij} \gamma_j^x) \rightarrow \sum_j (0, y_{ij} \gamma_j^x, z_{ij} \gamma_j^x) \quad (4.6)$$

$$\sum_j \mathbf{r}_{ij} \cdot \boldsymbol{\gamma}_j^y = \sum_j (x_{ij} \gamma_j^y, y_{ij} \gamma_j^y, z_{ij} \gamma_j^y) \rightarrow \sum_j (x_{ij} \gamma_j^y, 0, z_{ij} \gamma_j^y) \quad (4.7)$$

$$\sum_j \mathbf{r}_{ij} \cdot \boldsymbol{\gamma}_j^z = \sum_j (x_{ij} \gamma_j^z, y_{ij} \gamma_j^z, z_{ij} \gamma_j^z) \rightarrow \sum_j (x_{ij} \gamma_j^z, y_{ij} \gamma_j^z, 0) \quad (4.8)$$

Mathematical Formulations(contd.)

$$\mathbf{u}_i = -\frac{1}{4\pi} \sum_j \frac{1}{\sigma_j^3} g_1(w) (\mathbf{r}_{ij} \times \boldsymbol{\gamma}_j) \quad (4.9)$$

$$\mathbf{stx} = -\frac{1}{4\pi} \sum_j g_1(w) (\gamma_i^y \gamma_j^z - \gamma_i^z \gamma_j^y, \gamma_i^z \gamma_j^x - \gamma_i^x \gamma_j^z, \gamma_i^x \gamma_j^y - \gamma_i^y \gamma_j^x) \frac{1}{\sigma_j^3} \quad (4.11)$$

$$\mathbf{tx} = \frac{3}{4\pi} \sum_j g_2(w) (\boldsymbol{\gamma}_i \cdot \mathbf{r}_{ij}) (y_{ij} \gamma_j^z - z_{ij} \gamma_j^y, z_{ij} \gamma_j^x - x_{ij} \gamma_j^z, x_{ij} \gamma_j^y - y_{ij} \gamma_j^x) \frac{1}{\sigma_j^5} \quad (4.12)$$

$$\mathbf{I}_i = (\boldsymbol{\gamma}_i \cdot \mathbf{r}_i) \mathbf{S} - (\gamma_i^x \mathbf{T1} + \gamma_i^y \mathbf{T2} + \gamma_i^z \mathbf{T3}) \quad (4.14)$$

Function Table and Coefficients

Here,

$$w = \left(|\mathbf{r}_{ij}| / \sigma_j \right)^2; \quad |\mathbf{r}_{ij}| = |\mathbf{r}_i - \mathbf{r}_j|;$$

$g(w)$	A_j	B_j	\mathcal{E}_{ij}
$g1(w) = \frac{w+5/2}{(w+1)^{5/2}}$	$\frac{1}{\sigma_j^2}$	$\frac{\gamma_j}{\sigma_j^3}$	0
$g2(w) = \frac{w+7/2}{(w+1)^{7/2}}$	$\frac{1}{\sigma_j^2}$	$\frac{\gamma_j}{\sigma_j^5}$	0

Function Table

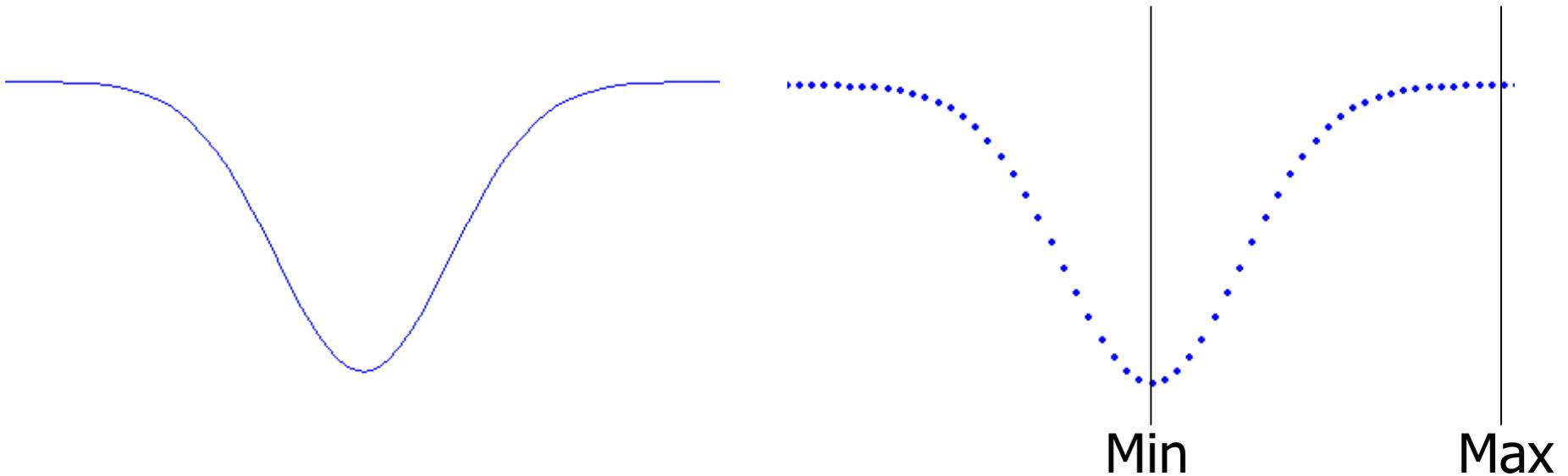
Function $g(w)$:

$g(w): \min \leq w \leq \max$ (Domain of function)

$$\frac{w_{\max}}{w_{\min}} \leq 2^{32}$$

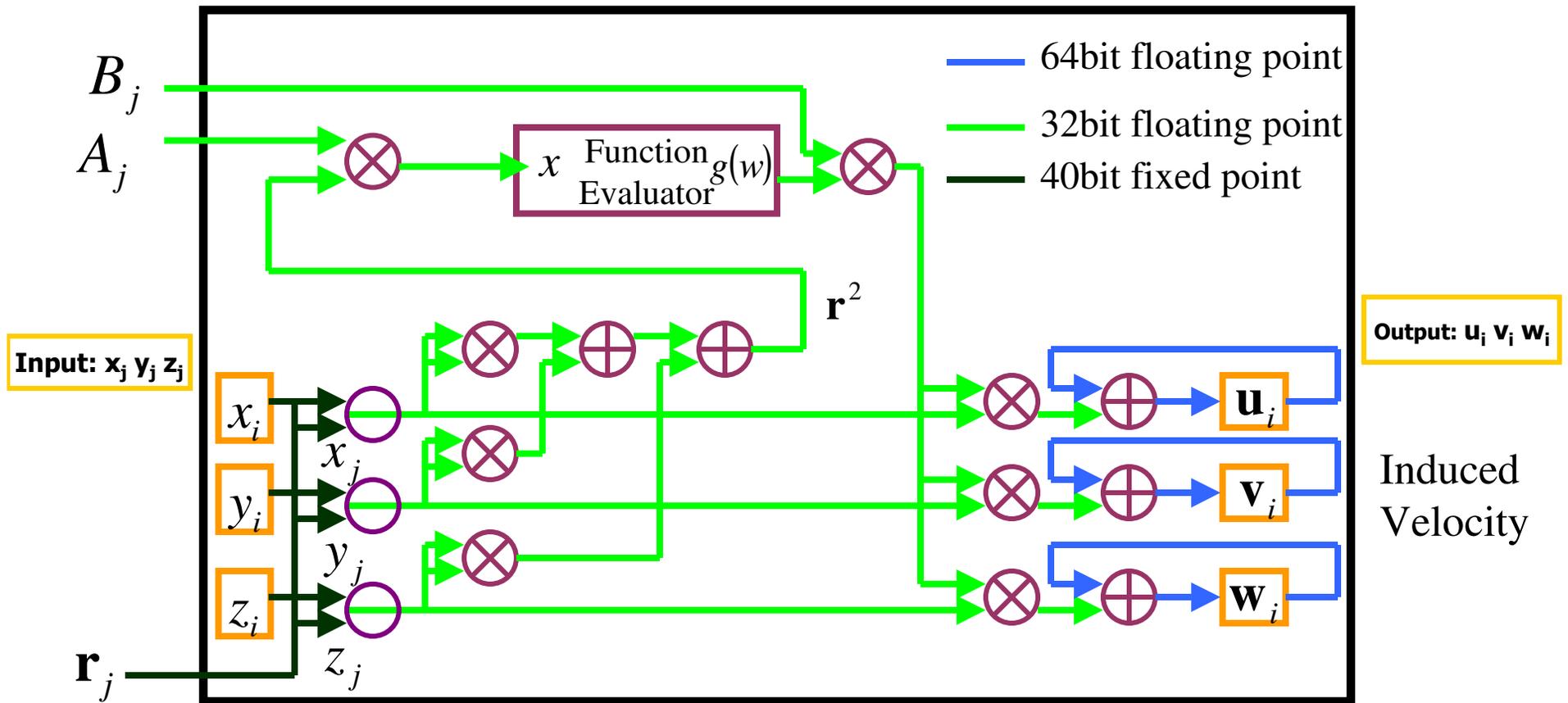
$g(w)$

Function Table



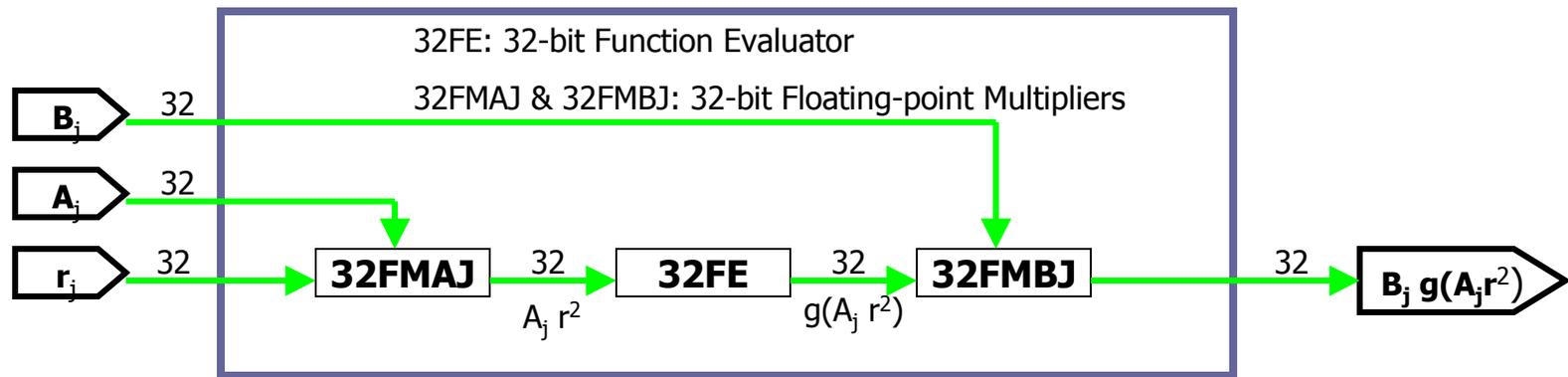
Block Diagram of MDGRAPE-2 pipeline

$$\text{Pairwise force: } \vec{f}_{i,j} = b_{ij} g(a_{ij} r_{ij}^2) \vec{r}_{ij} \quad (3.1)$$



Ref: Tetsu Narumi, 1997, PhD Thesis, Tokyo University

Function Evaluator



Approximation:

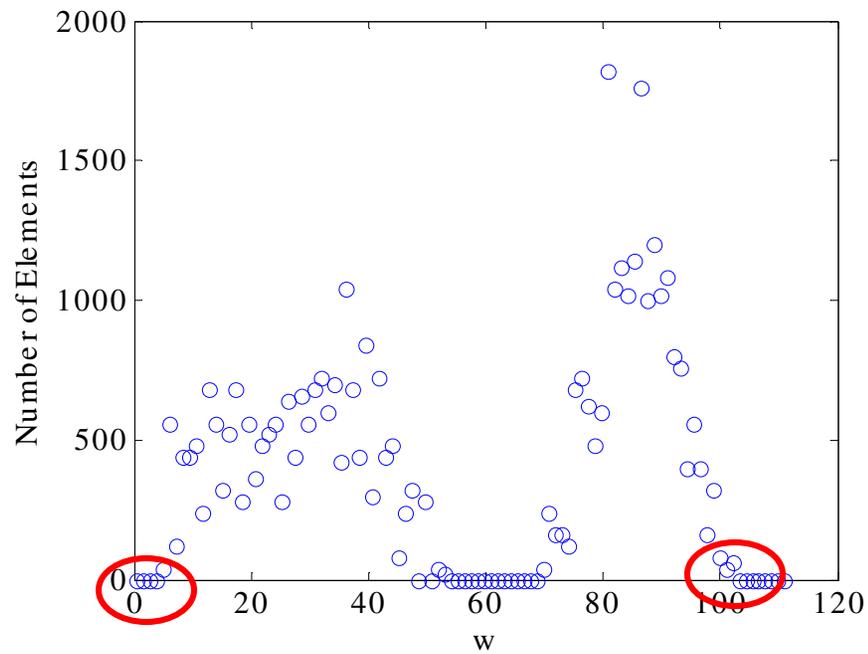
$$g(w) = c_0 + w(c_1 + w(c_2 + w(c_3 + wc_4)))$$

➡ 32 bit floating-point calculation

➡ Relative accuracy: 10^{-7}

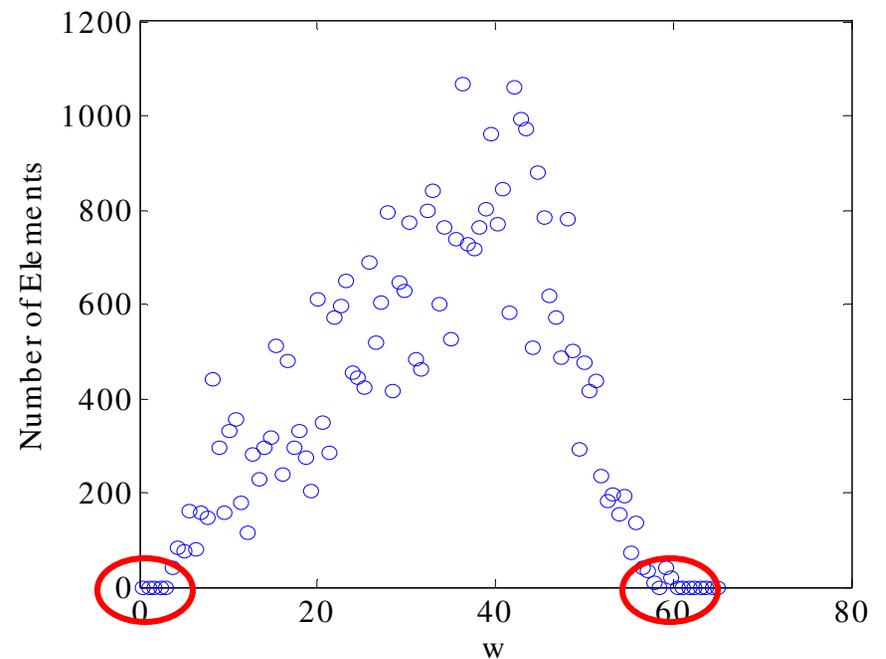
4.3 Typical distribution of vortex elements

$$t\Gamma/R^2 = 1$$



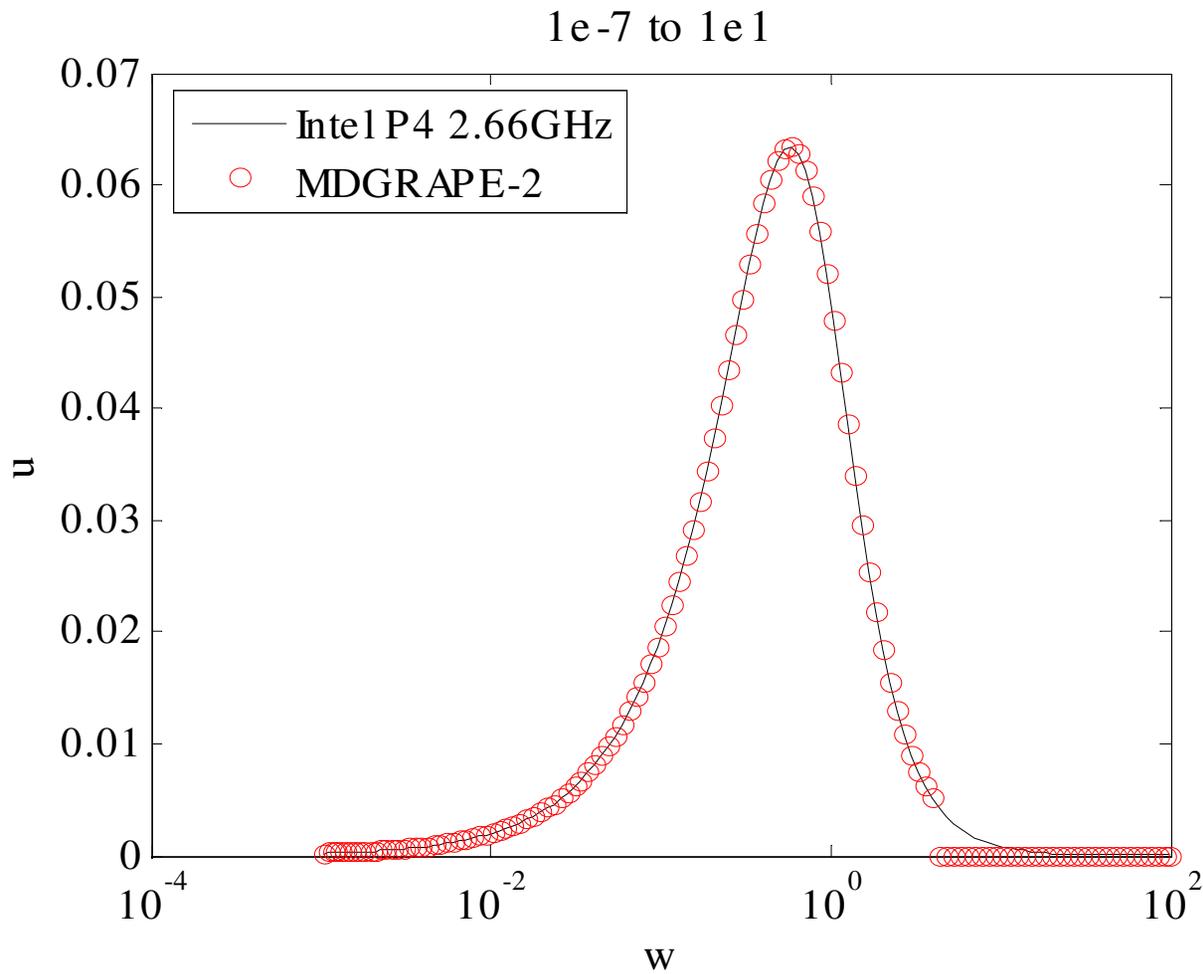
Min:3.9; Max: 103.2

$$t\Gamma/R^2 = 100$$



Min:2.94; Max: 60.37

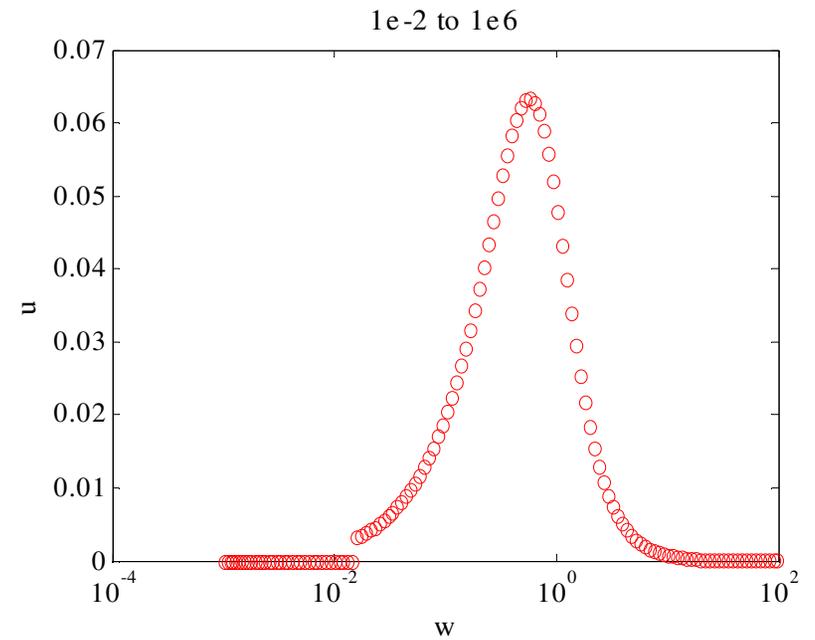
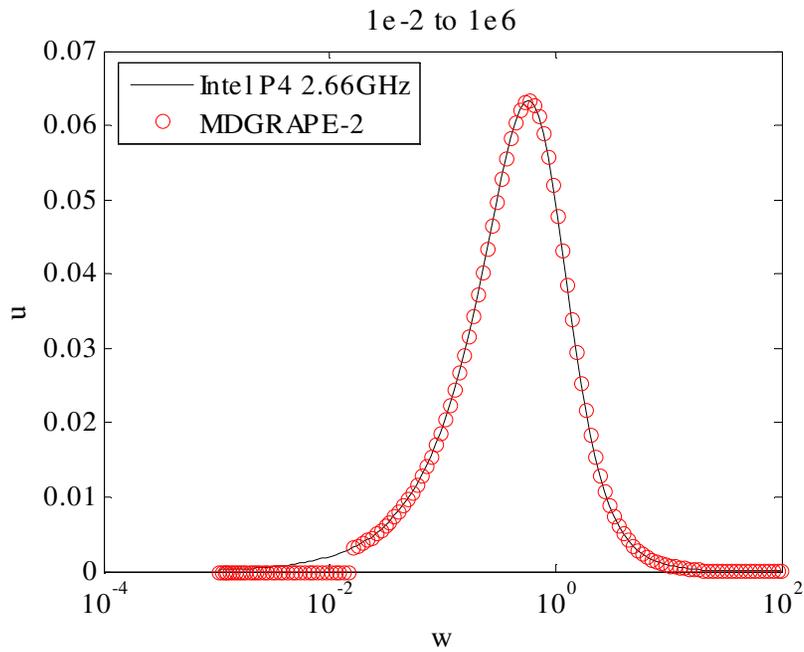
Range of a Function Table



tested table ranges

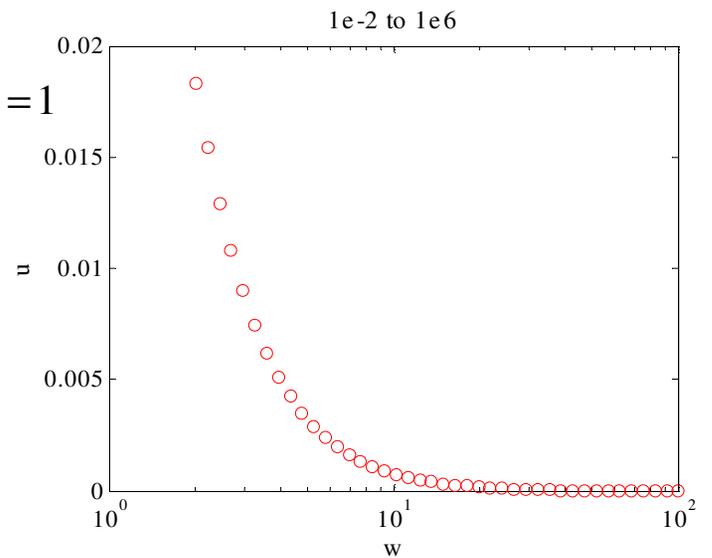
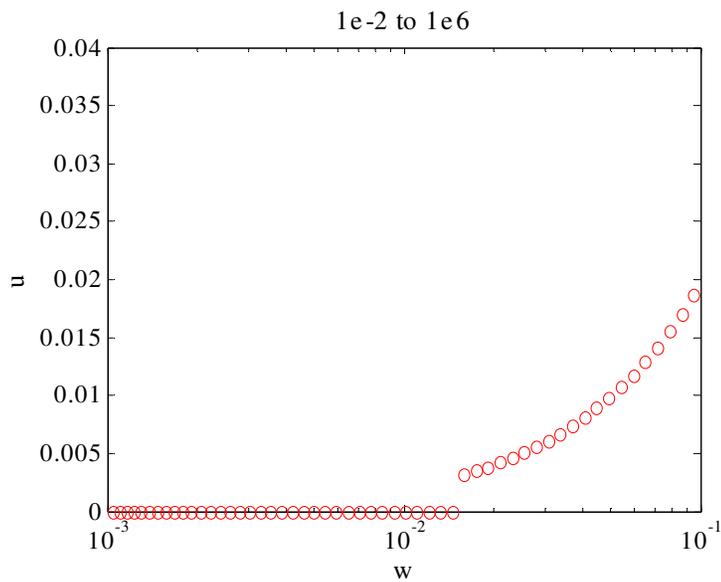
min	max
10^{-2}	10^6
10^{-3}	10^5
10^{-4}	10^4
10^{-5}	10^3
10^{-6}	10^2
10^{-7}	10^1

4.3.4 Optimum Range of a Function Table



$$2^{-12} < w < 2^{20}$$

$$w = (r_{ij} / \sigma_j)^2; \text{ here } \sigma = 1$$



4.5 Application

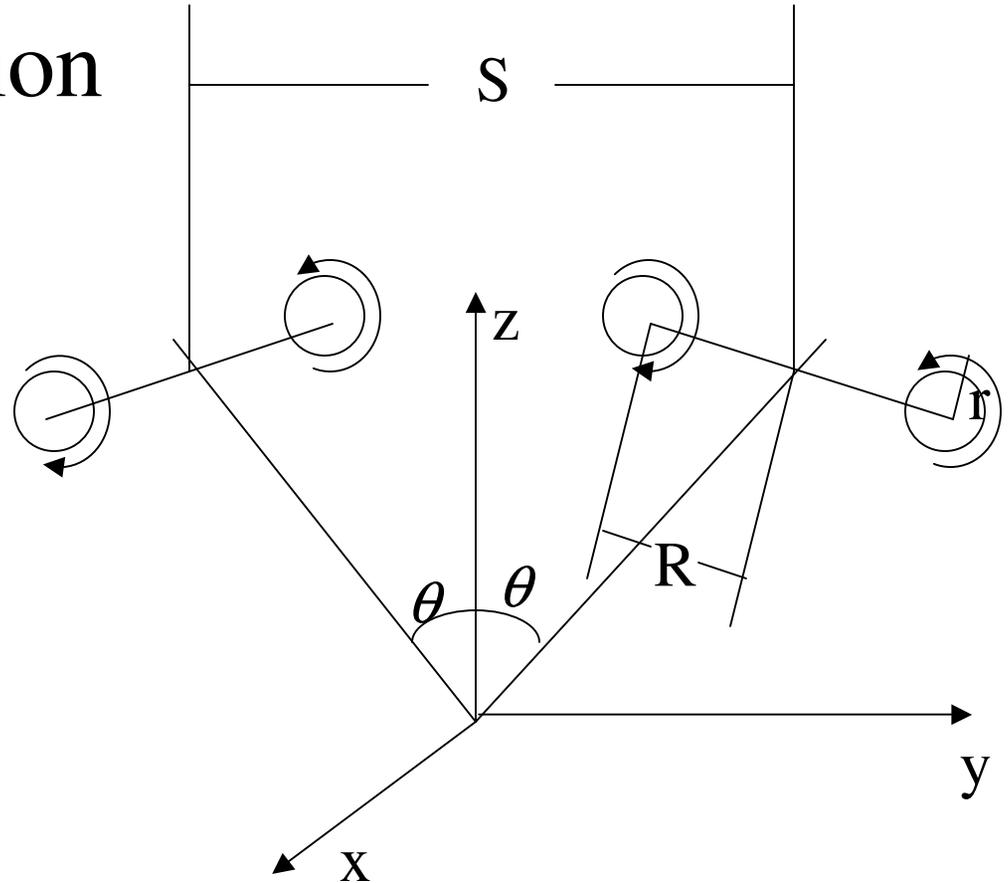
Calculation Condition

R = Radius of ring

r = Radius of cross-section

S = Distance between the center
of two rings

θ = Inclined angle of a ring

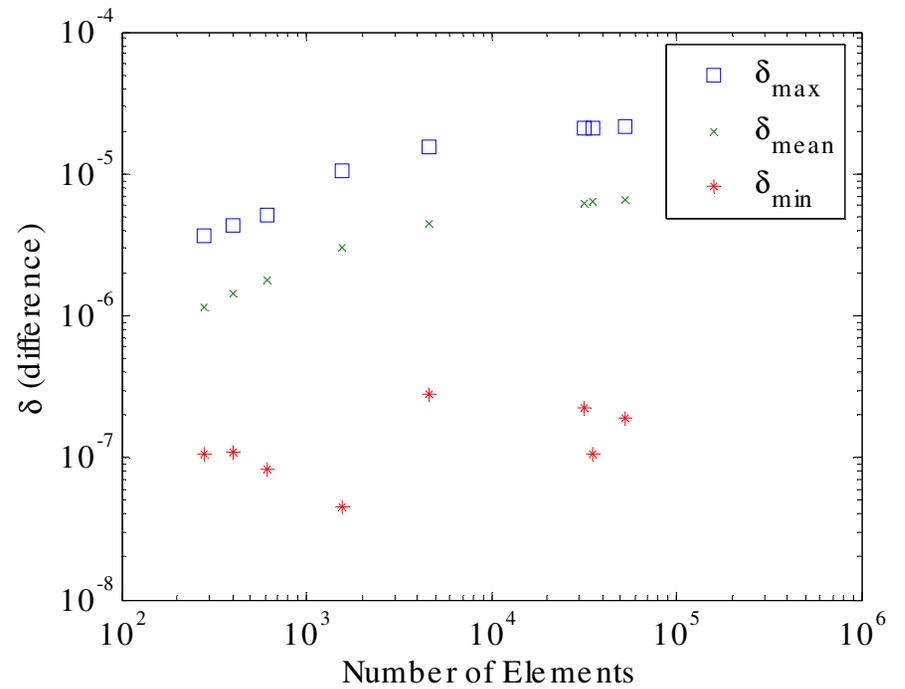
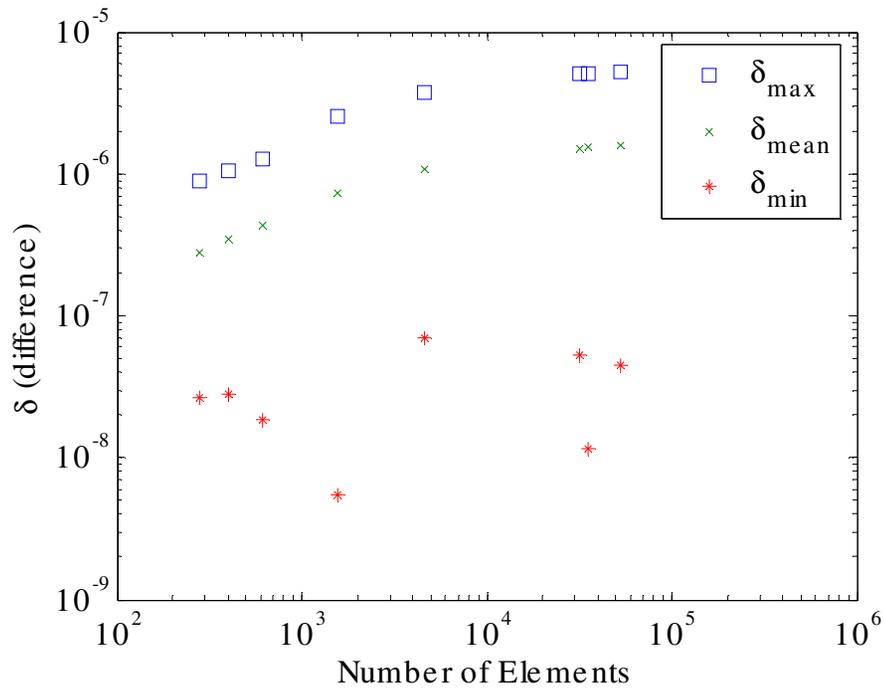
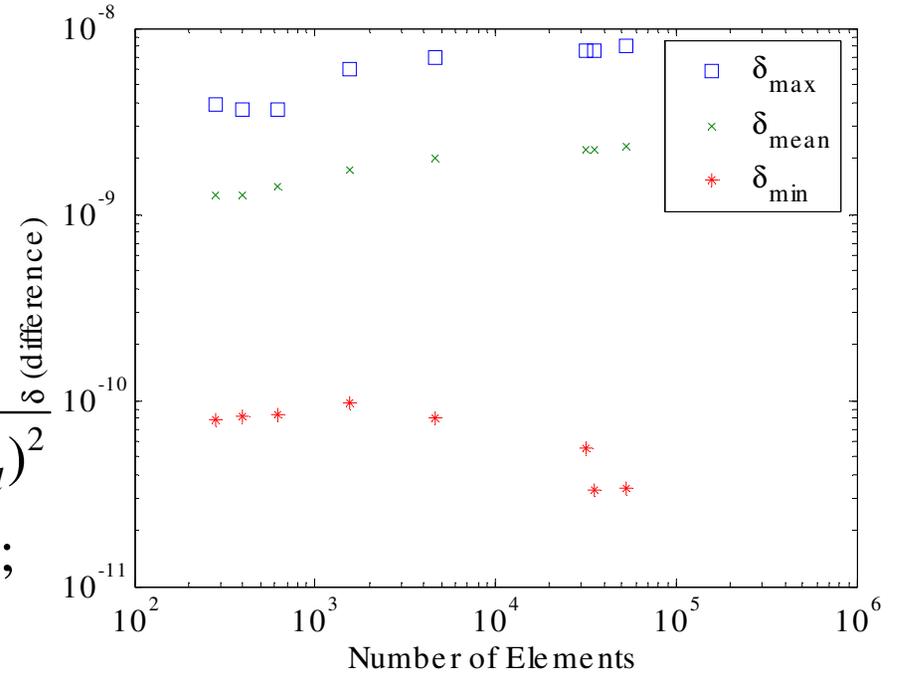


Position	R	r	σ	Γ_0	N	Re_{Γ}	S	θ
Inclined	1	0.05	0.065	1	$2 \times 502 \times 61$	400	2.7	15°

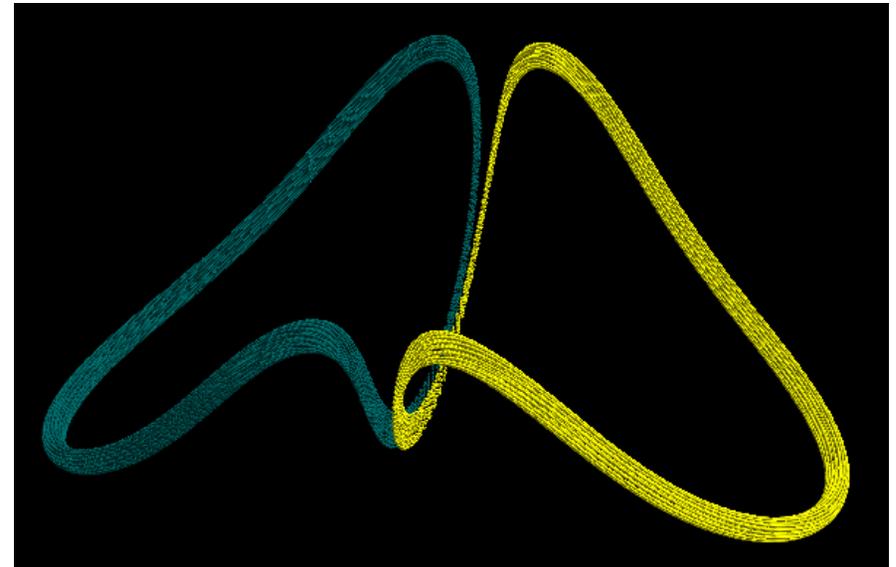
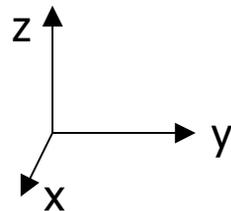
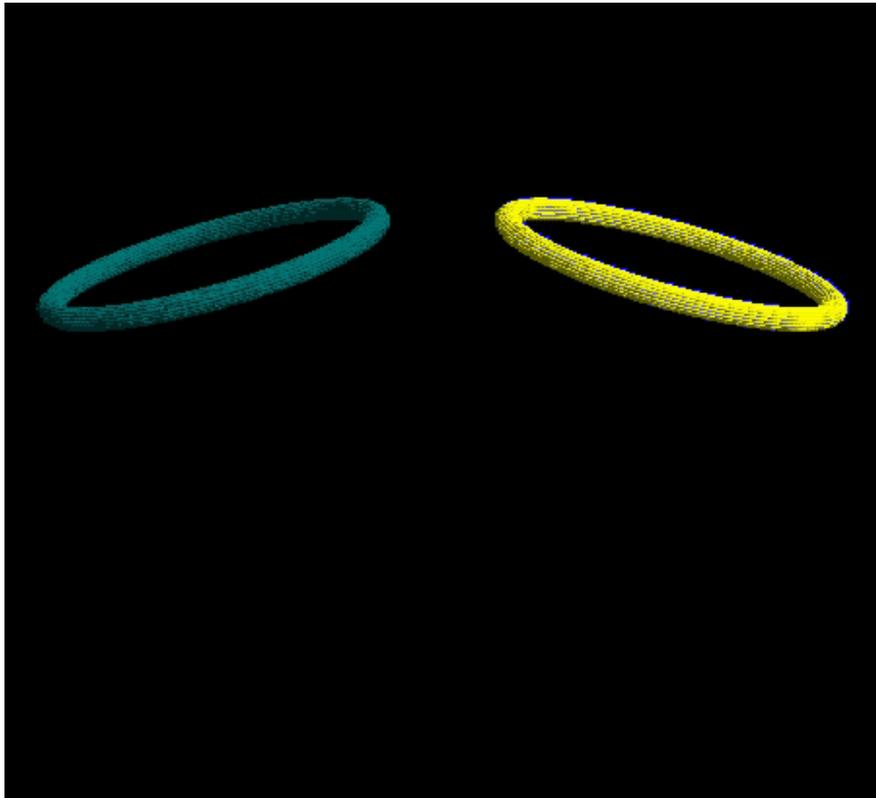
Convection error

$$\delta = \sqrt{(x_{host} - x_{md})^2 + (y_{host} - y_{md})^2 + (z_{host} - z_{md})^2}$$

$$\delta_{\min} = \min(\delta); \delta_{\text{mean}} = \text{mean}(\delta); \delta_{\max} = \max(\delta);$$



Snapshots of vortex elements



Ring radius $R=1$

Cross-section radius $r=0.05$

Reynolds number $Re=400$

Core radius = 0.065, Circulation of ring = 1

Total number of elements = $2 \times 502 \times 61$

Particles are evenly distributed

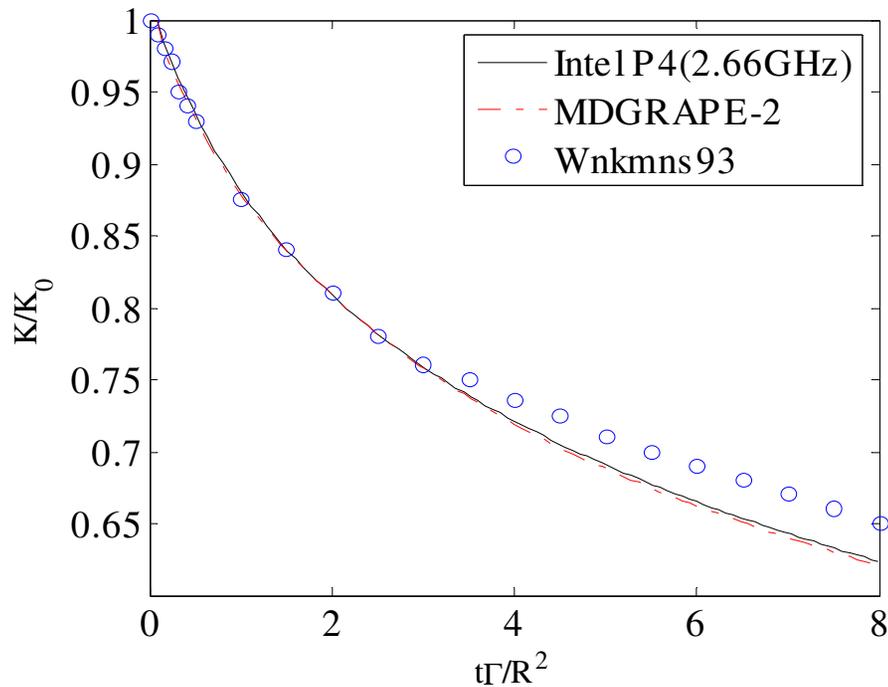
Initial distance between two rings = 2.7

Inclined angle = 15°

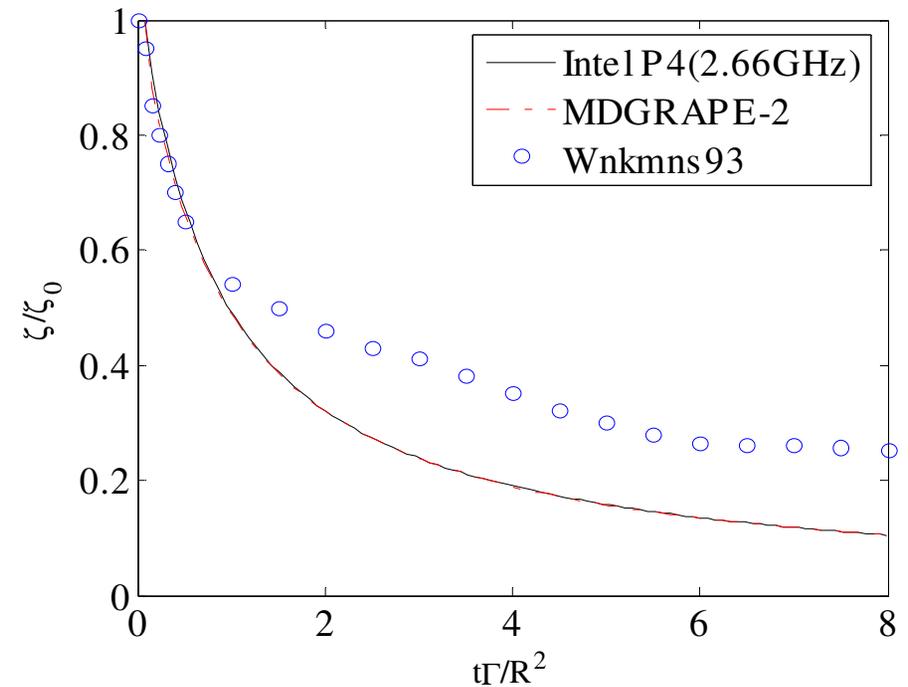
Kinetic Energy & Enstrophy

Kinetic energy: Eq. (4.16) and Enstrophy: Eq. (4.17)

Kinetic Energy



Enstrophy



Wnkmns93: G. S. Winckelmans and Leonard, J. Comp. Phys, 109, 247-273(1993)

Summary

- A mathematical formulation has been developed using MDGRAPE-2
- A rigorous assessment of this hardware has been made for a pair of impinging vortex rings
- Computational domain has been investigated that determines optimal range of a function table
- The global kinetic energy and enstrophy has been evaluated to address the numerical accuracy
- The results have good agreement when compared with the host calculation and referenced work

5

The Study of Colliding Vortex
Rings using a Special-purpose
Computer and FMM

5.1 Introduction

- Simultaneous use of the FMM with MDGRAPE-3
- To investigate the possibility of further accelerations
- The various forms of FMM are investigated
- The accuracy is achieved by simulating the impingement of two identical inclined vortex rings
- The effect of temporal and spatial resolutions will be investigated
- The reconnection of vortex rings is observed

Special-purpose Computer: MDGRAPE-3

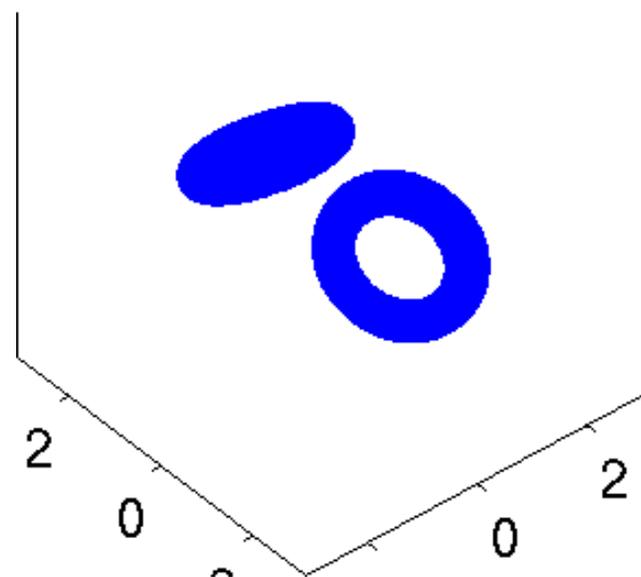
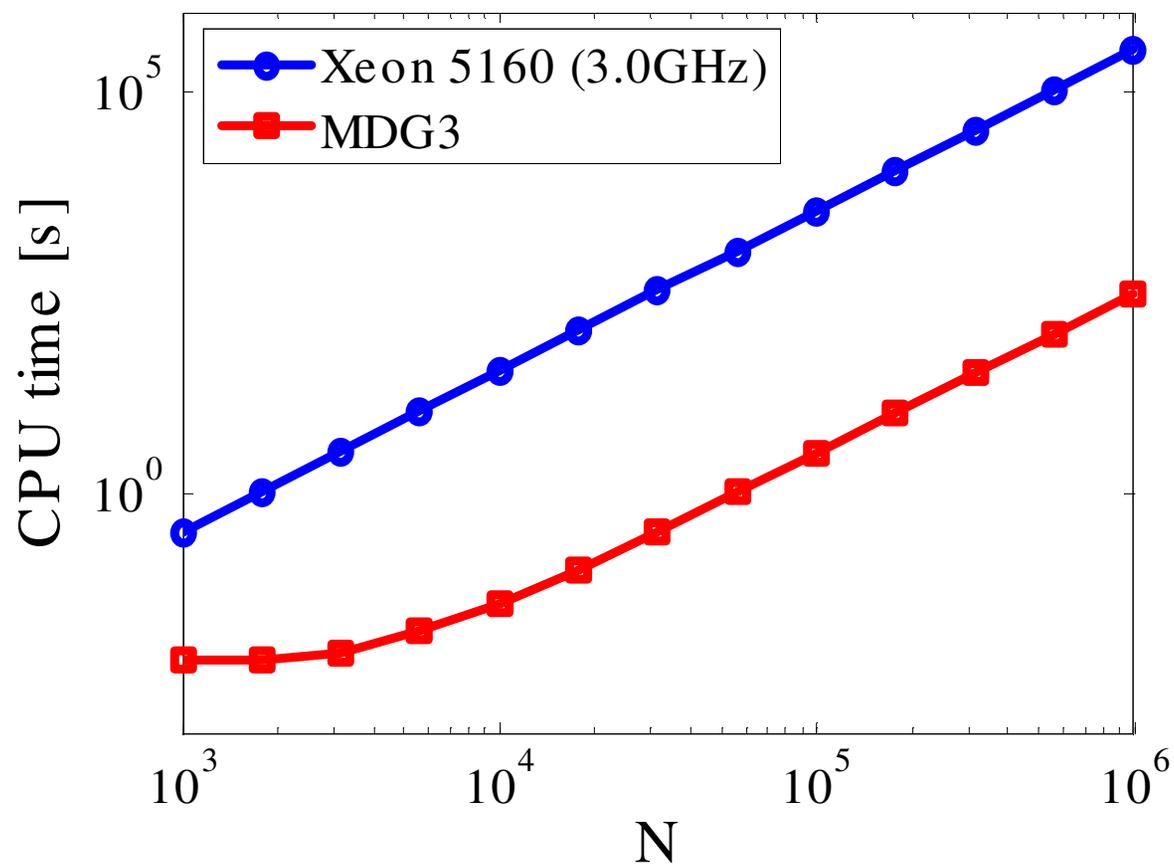
(Taiji, 2003; Narumi, 2006)

- Petaflops special-purpose computer and the successor of MDGRAPE-2
- One small board consists of 2 chips
- One MDGRAPE-3 chip combined with 20 parallel pipelines
- Calculation speed of one small board is 330GFlops/250MHz
- 12.5 times faster than MDGRAPE-2
- Total memory/chip: 9Mbits

MDGRAPE-3 Board



Performance



$N=10^6$

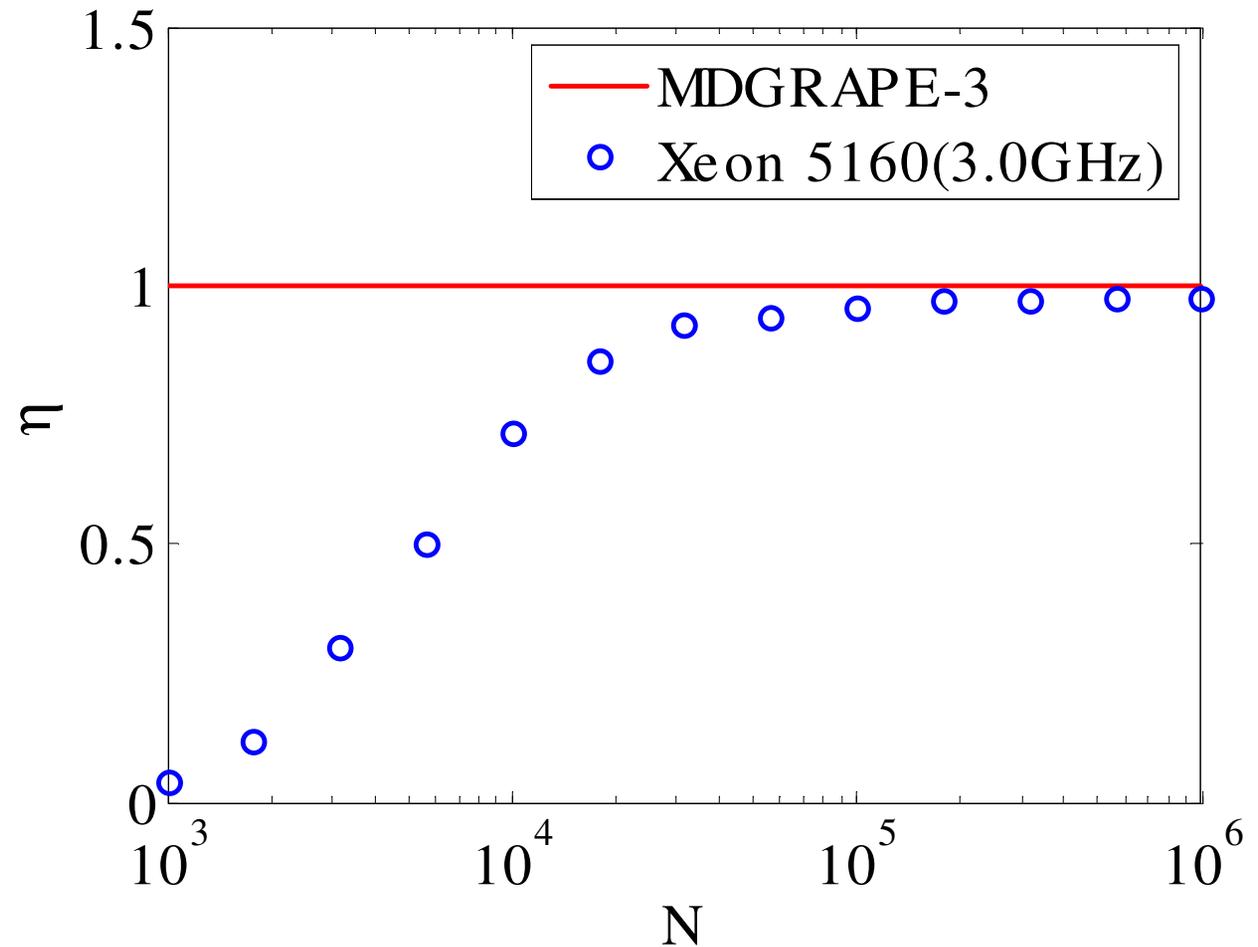
1000 times at $N \sim 10^6$

Efficiency of VM Calculation

$$N_{APPL} = \frac{nmd(xmd + ymd + zmd)}{CPUtime(sec/step)}; \quad N_{GRAPE} = 10^{10}$$

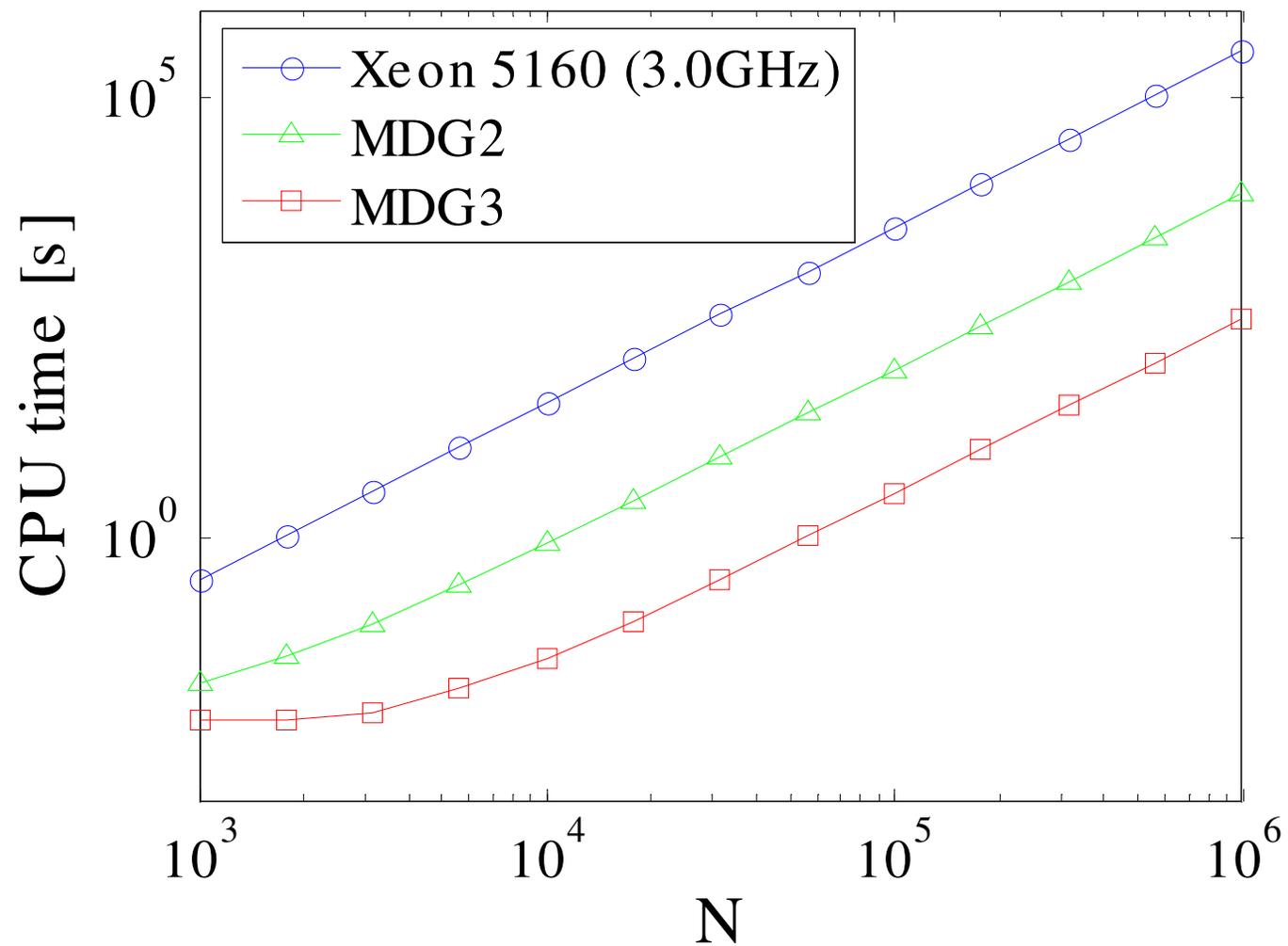
Efficiency:

$$\eta = \frac{N_{APPL}}{N_{GRAPE}}$$

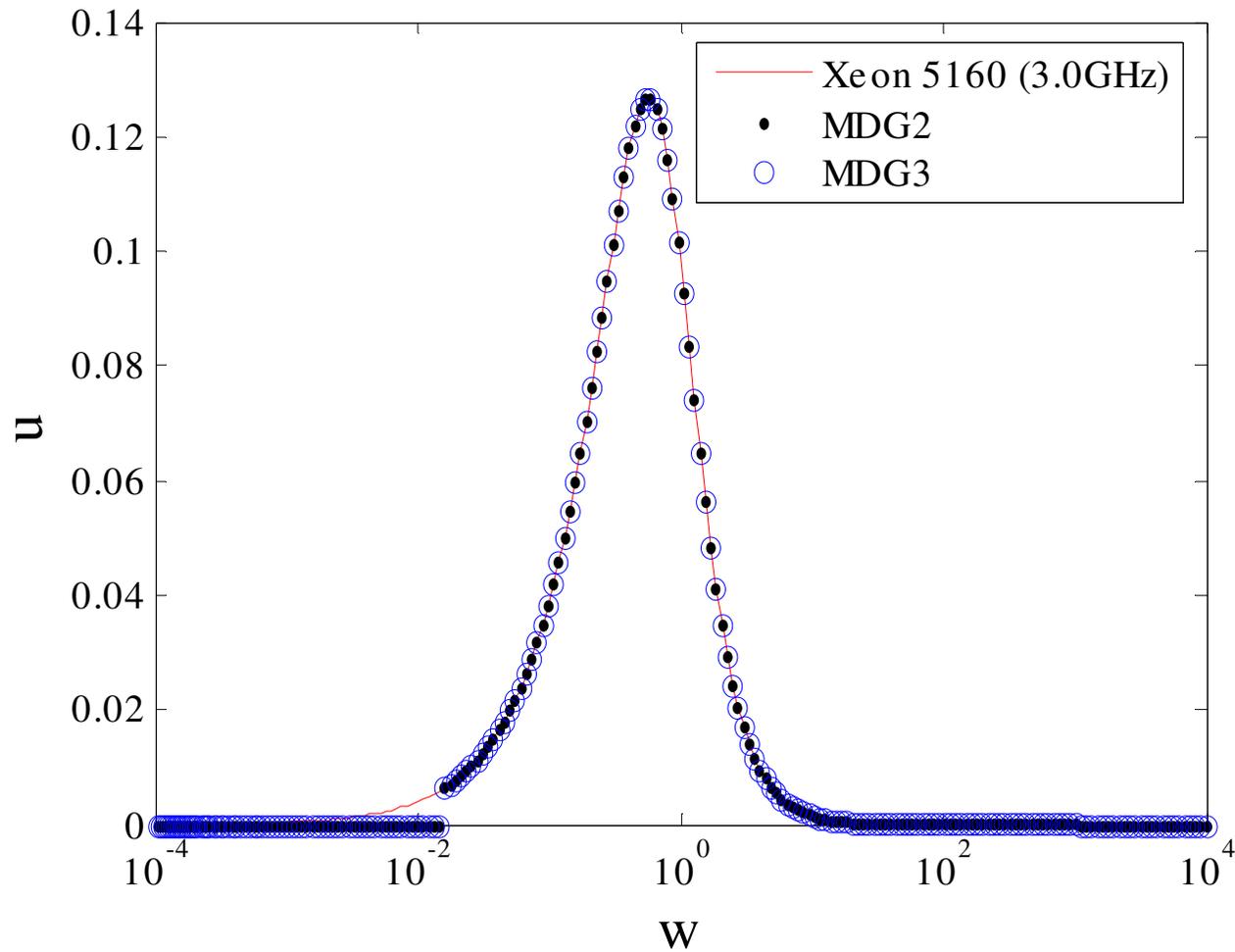


Comparative study between
MDGRAPE-2 and MDGRAPE-3

CPU-Time

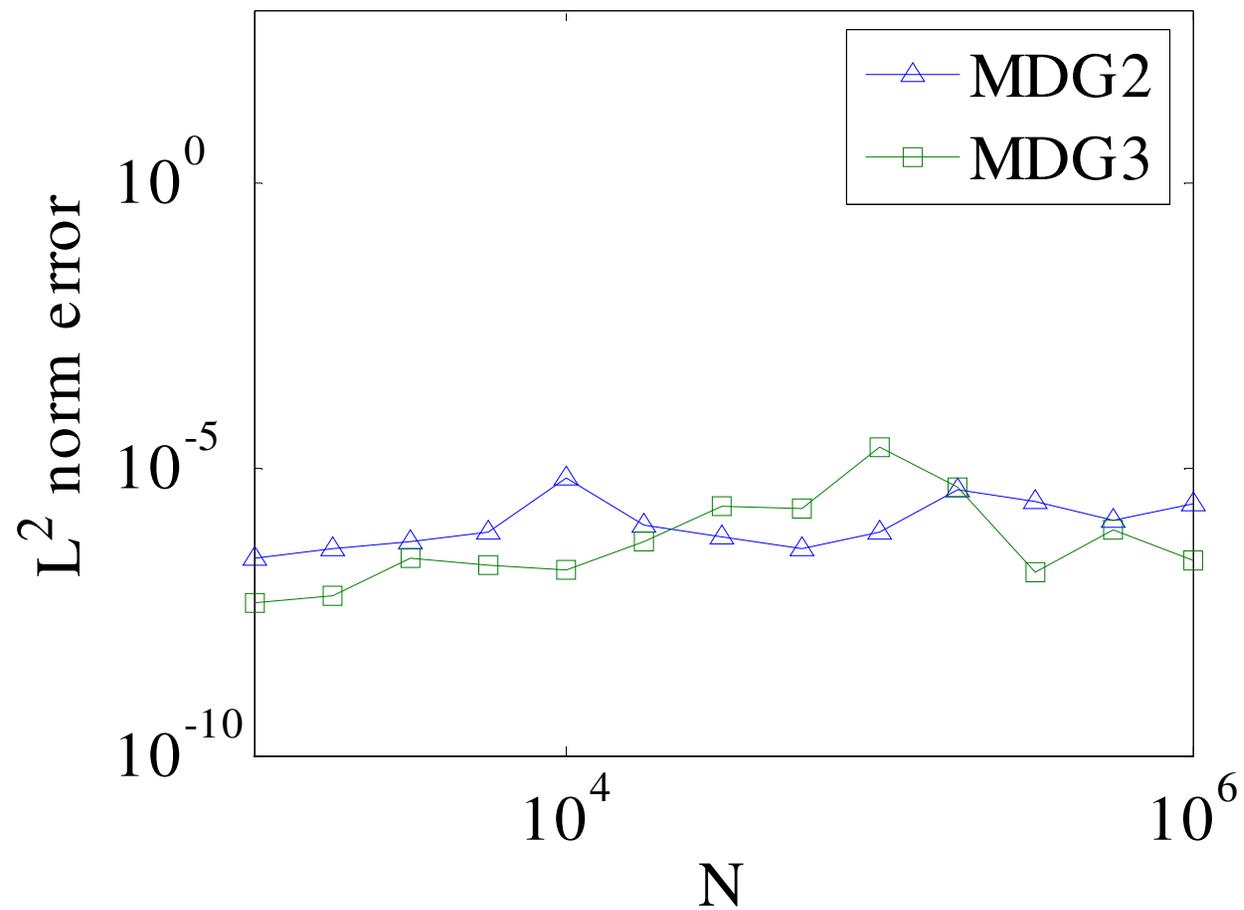


Scaling Error



L^2 norm error

$$L^2 = \frac{\sum_j \left\{ (u_{host} - u_{md})^2 + (v_{host} - v_{md})^2 + (w_{host} - w_{md})^2 \right\}}{\sum_j \left\{ u_{host}^2 + v_{host}^2 + w_{host}^2 \right\}} \quad (3.14)$$



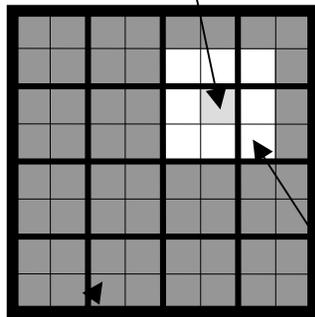
FMM on MDGRAPE-3

Hot-Spot of FMM

red is source → blue is target

General rule : M2L can not be preformed for neighbors

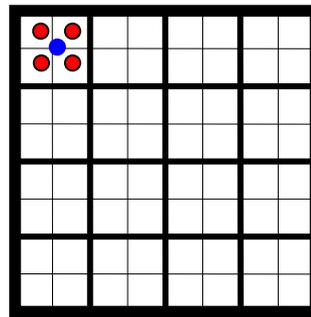
If we want to calculate for this box



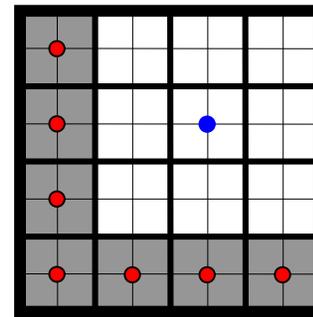
Neighbor particles are solved directly

Far particles are solved by the FMM

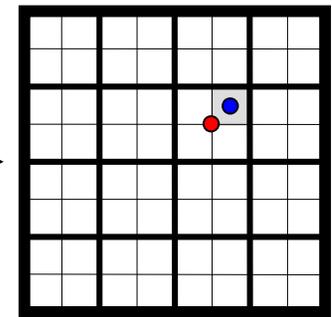
M2M



M2L

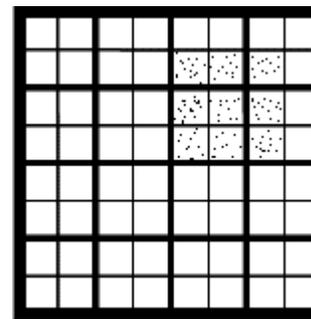


L2L

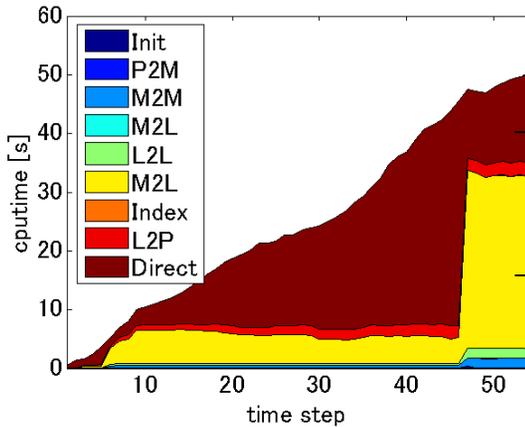
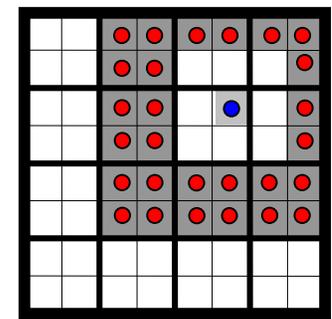


+

Direct



M2L



Direct

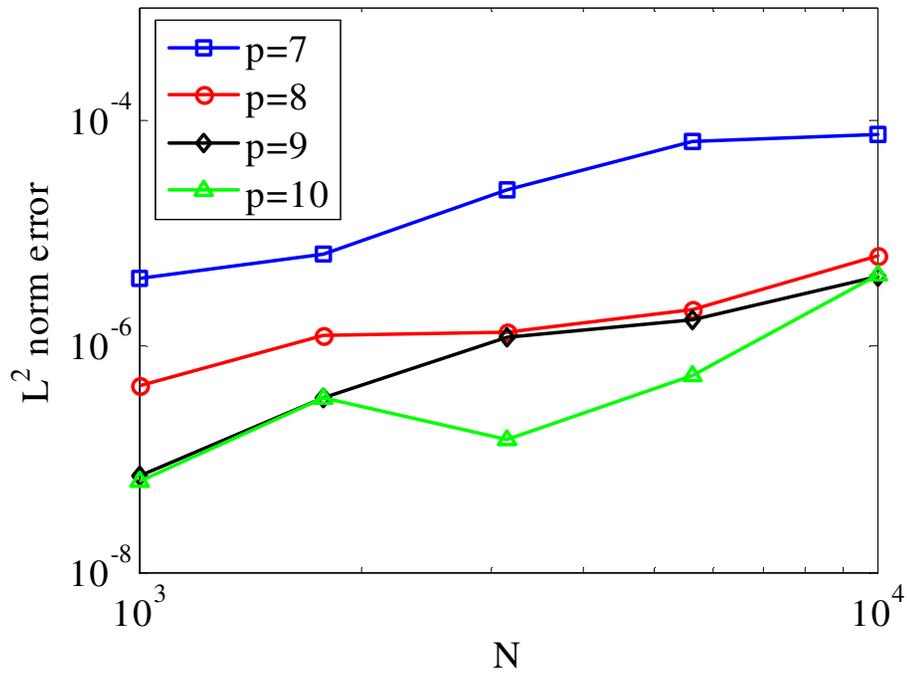
M2L

If the box is too coarse this could also be the hot-spot

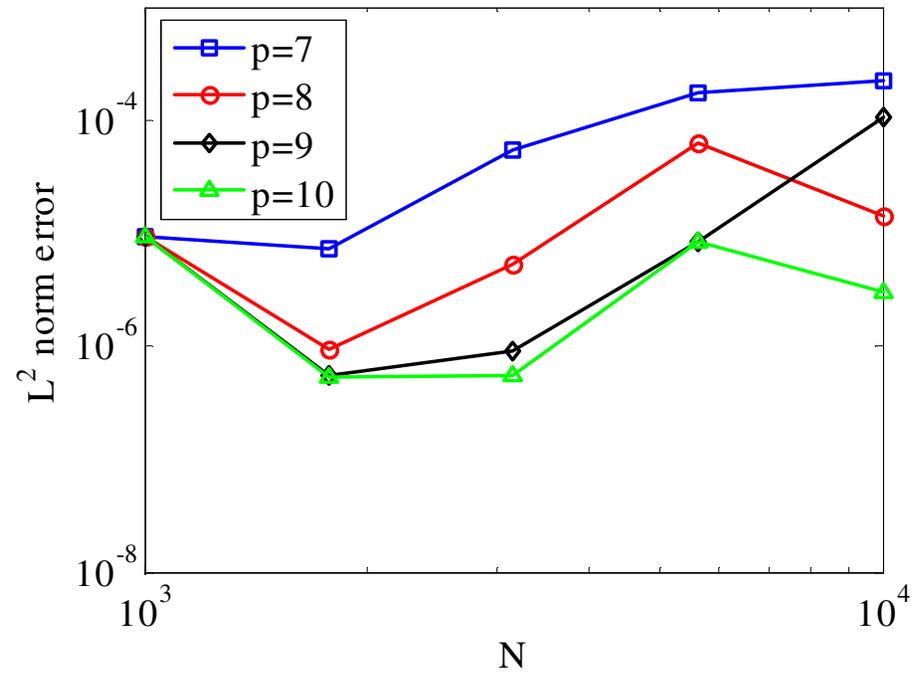
Many sources acting on one target
This is the hot-spot of the FMM

Momentum Effect on FMM Accuracy

$$\mathbf{u}_i \approx \frac{1}{4\pi} \sum_{n=0}^p \sum_{m=-n}^n \left\{ \sum_{j=1}^N \gamma_j M_j \right\} \times \nabla S_i \quad (2.13)$$



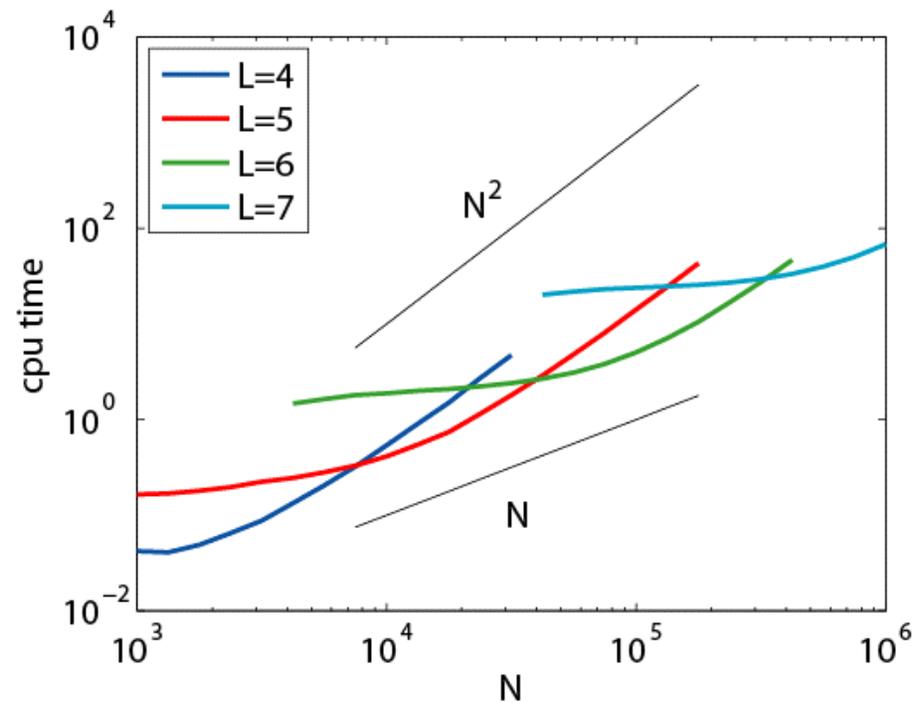
Biot-Savart



Stretching

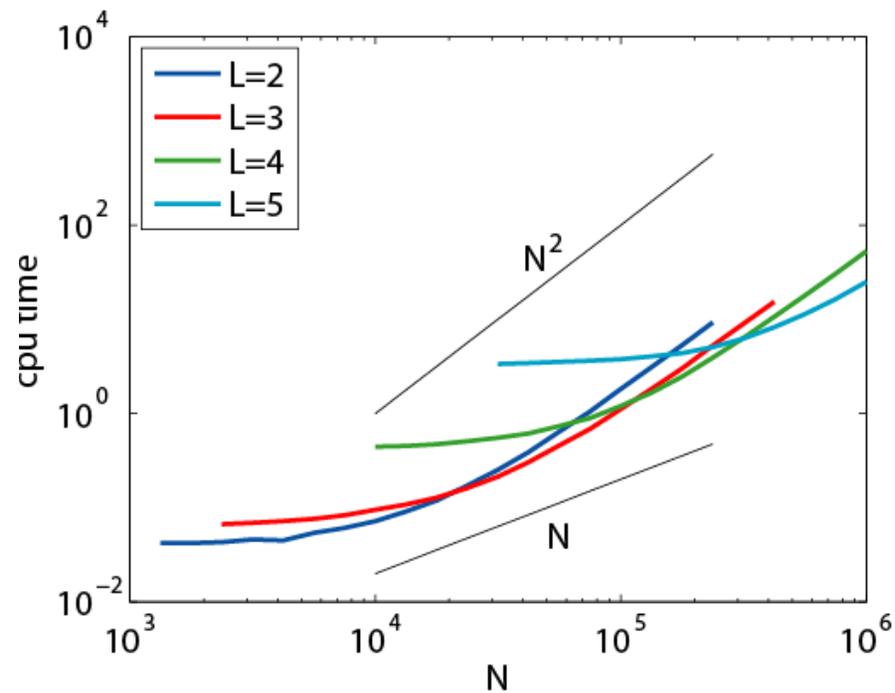
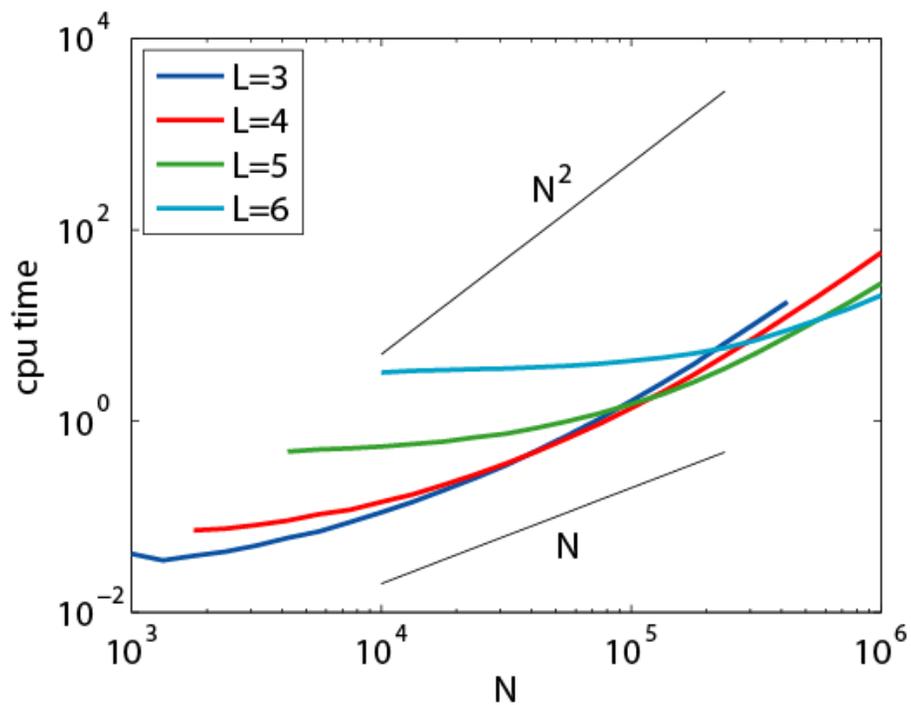
5.4.1 Optimization

FMM

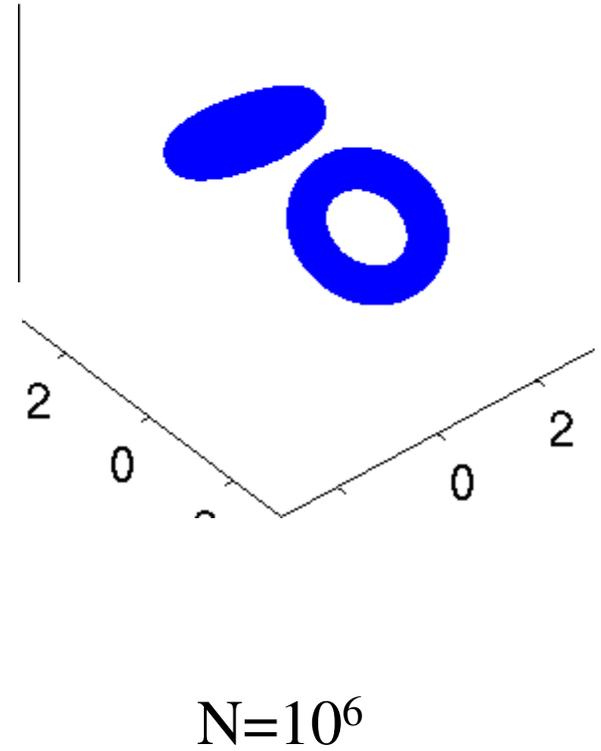
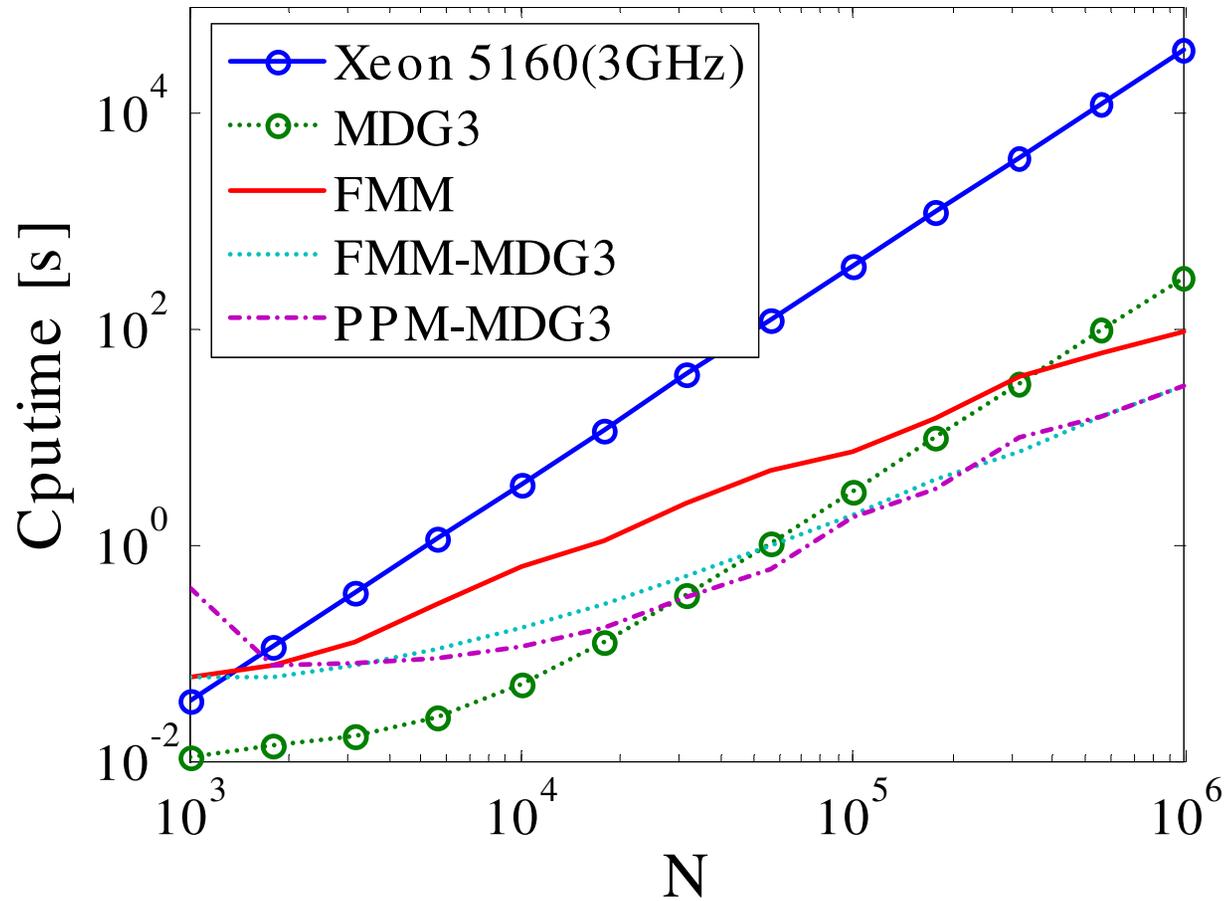


FMM-MDG3

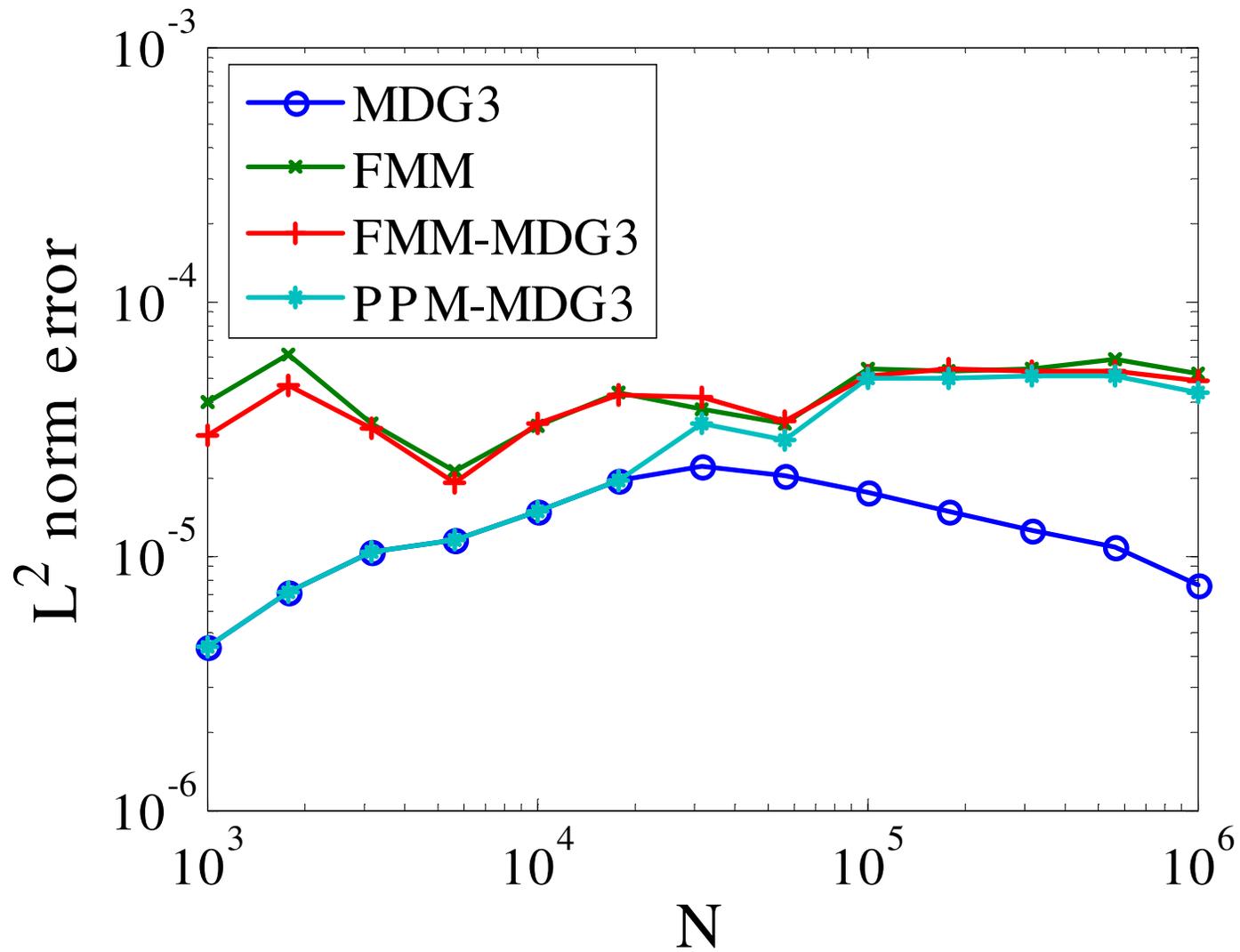
PPM-MDG3



5.4.2 CPU-Time



5.4.2 Error



Acceleration Ratio

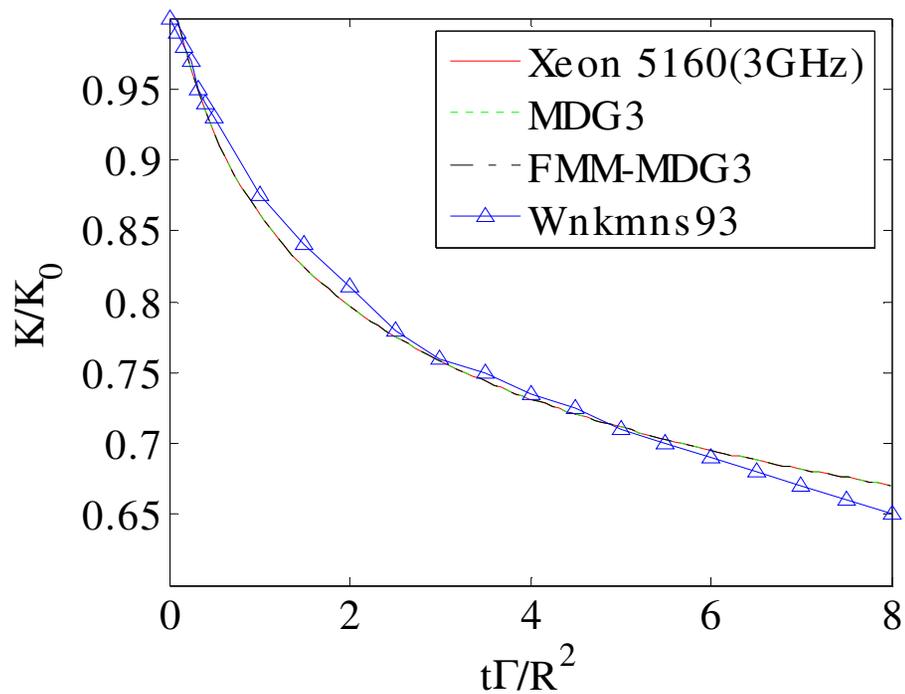
$N=10^6$

Biot-Savart		Stretching	
Direct		Direct	
↓ ×462	↓ ×119	↓ ×613	↓ ×52
FMM	MDG3	FMM	MDG3
↓ ×4.1	↓ ×16	↓ ×2.8	↓ ×33
FMM+MDG3		FMM+MDG3	

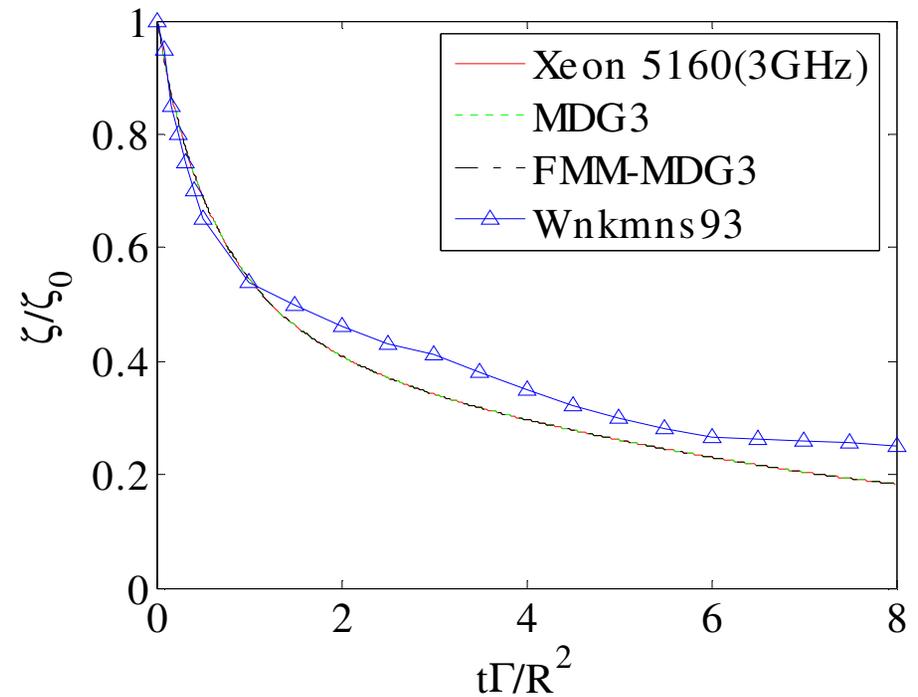
5.5 Vortex Ring Calculation

Energy and Enstrophy

Eq. (4.16)

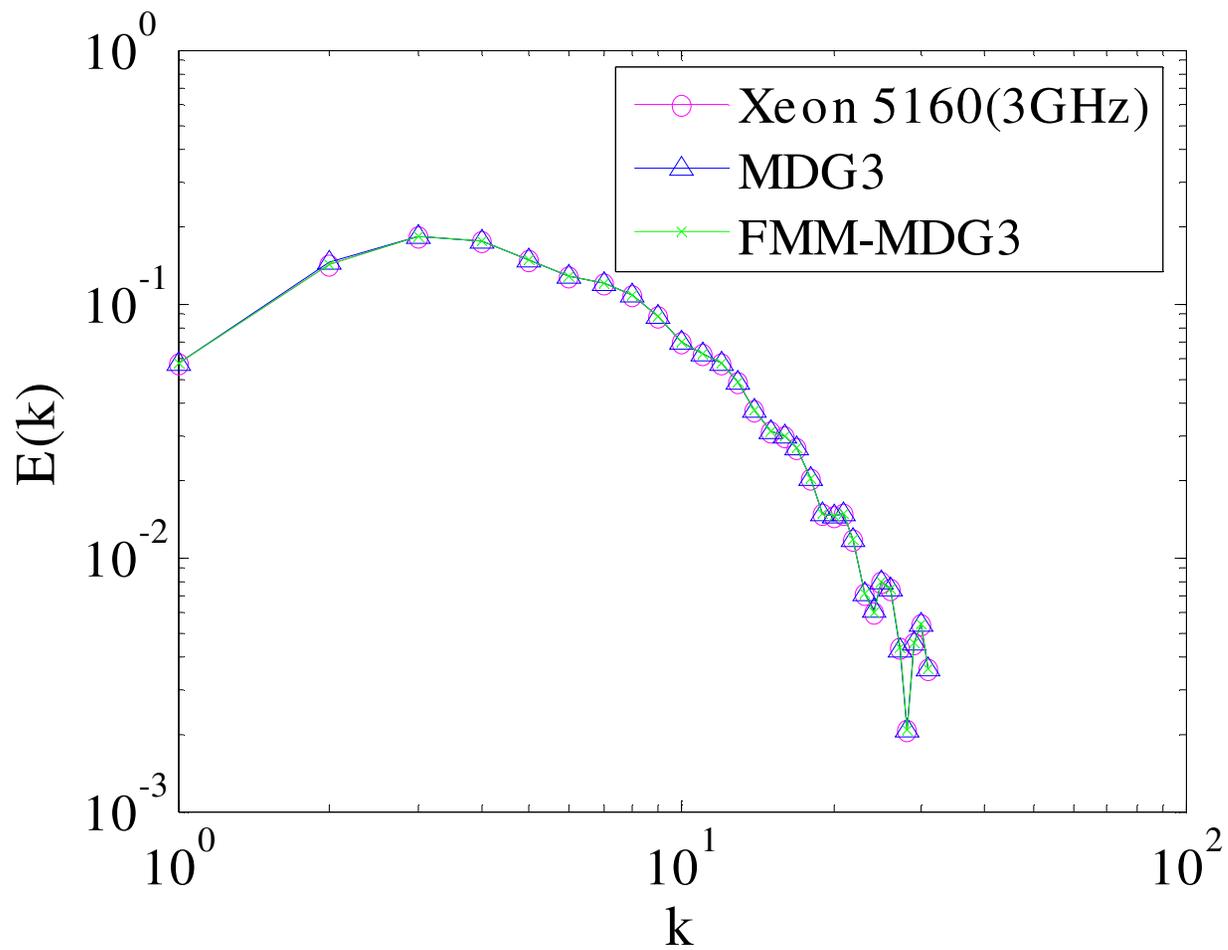


Eq. (4.17)



Wnkmns93: G. S. Winckelmans and Leonard, J. Comp. Phys, 109, 247-273(1993)

Energy Spectra



5.5.3 Calculation Conditions

- Vorticity distribution:
$$\boldsymbol{\omega} = \frac{\Gamma}{2\pi\sigma^2} \exp\left(\frac{-\mathbf{r}^2}{2\sigma^2}\right) \quad (5.2)$$

- Initial core radius:
$$\frac{\sigma_0}{h} = 2 \quad (5.3)$$

- Kinetic Energy:
$$K = \frac{1}{2} \sum_i^N \mathbf{u}_i \cdot \mathbf{u}_i \quad (5.4)$$

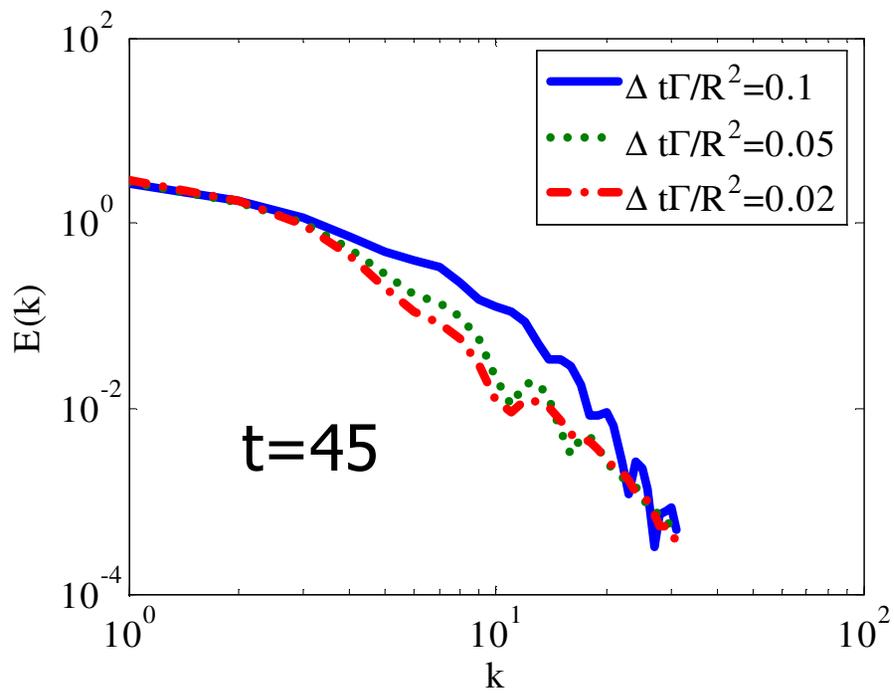
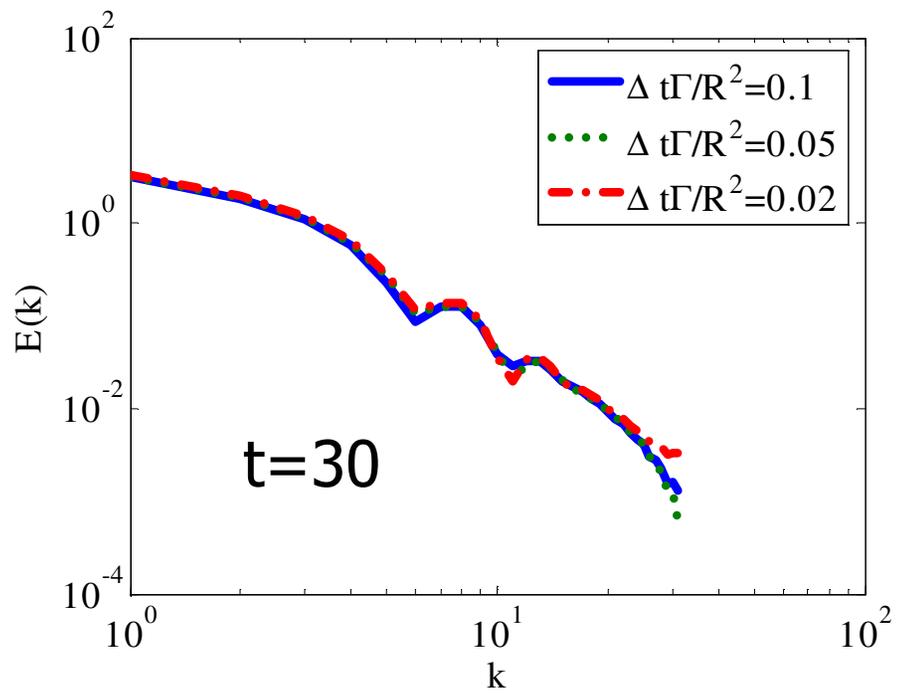
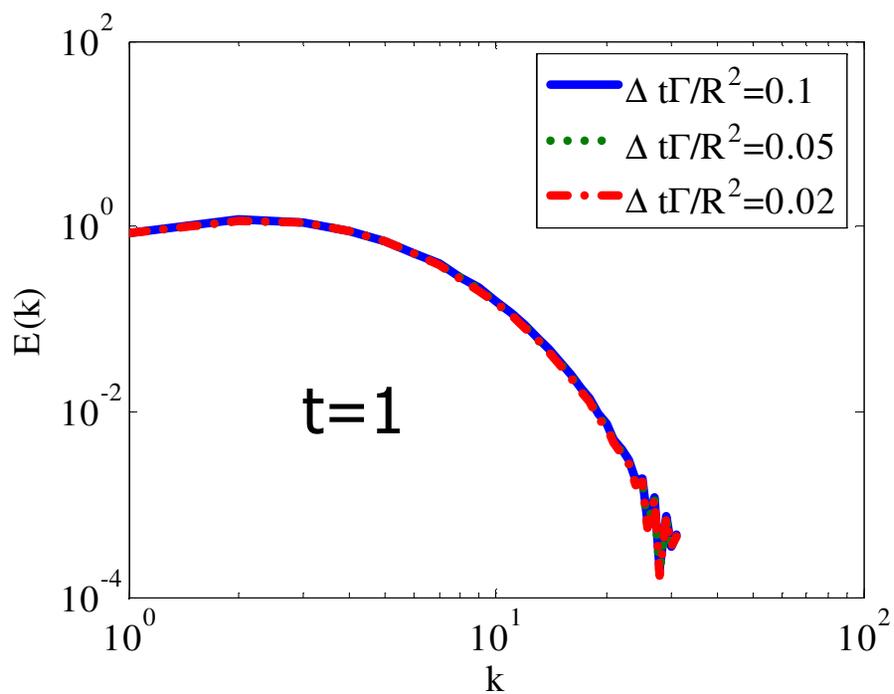
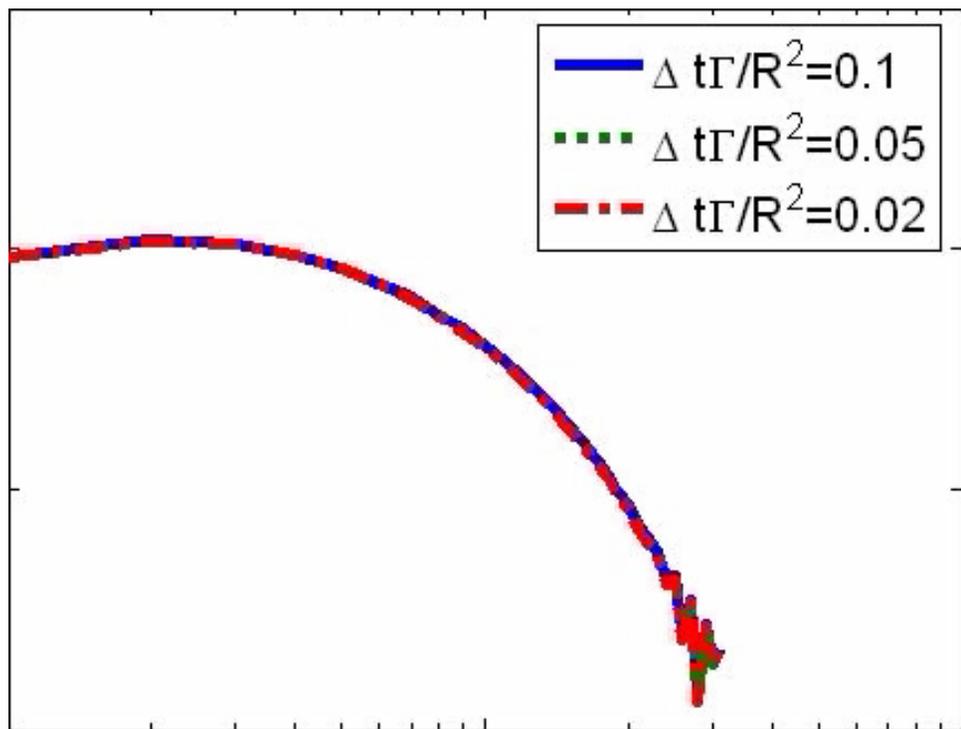
- Enstrophy:
$$\zeta = \sum_i^N \boldsymbol{\omega}_i \cdot \boldsymbol{\omega}_i \quad (5.5)$$

- Reynolds number:
$$\text{Re}_\Gamma = \frac{\Gamma}{\nu} = 400$$

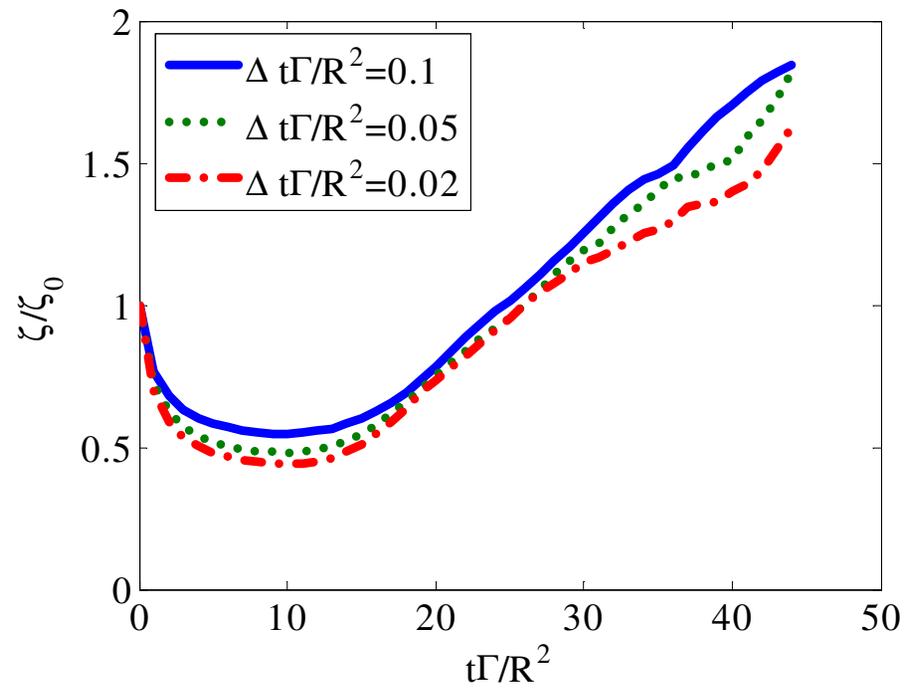
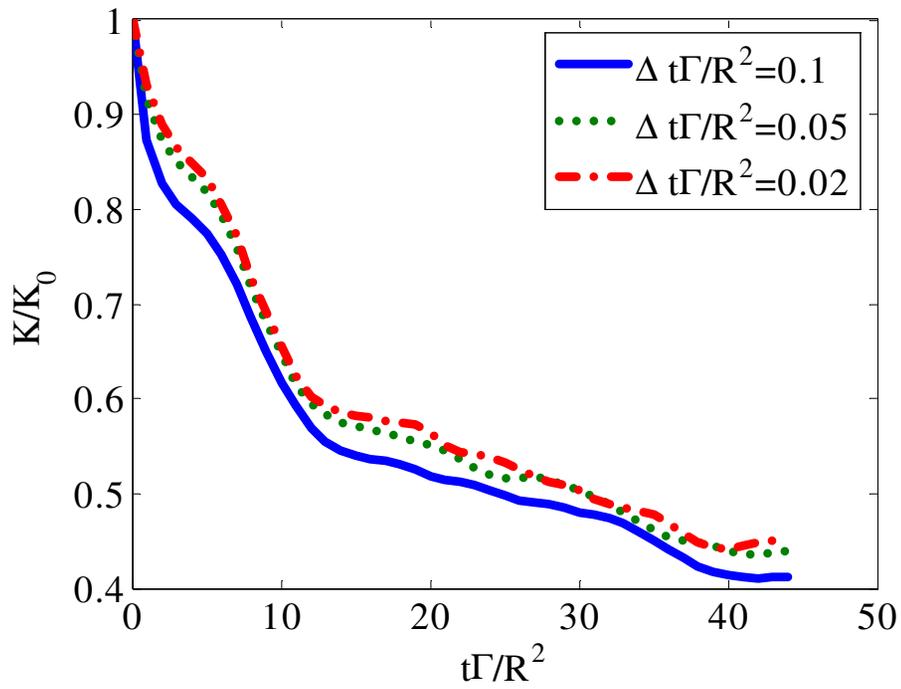
- Number of particles:
$$N = 10^5 \sim 10^7$$

5.5.4 Effect of Temporal Resolution

$$\frac{\Delta t \Gamma}{R^2} = 0.1, 0.05, 0.02$$



Kinetic Energy and Enstrophy (dt)



5.5.5 Effect of Spatial Resolution

Case	A	B	C
Number of Rings	2	2	2
N per Cross Section	190	418	910
Cross Sections	271	1261	5677
Total	102980	1054196	10332140

Position of vortex elements

Ring radius $R=1$; Cross-section radius $r=0.05$

Reynolds number $Re=400$

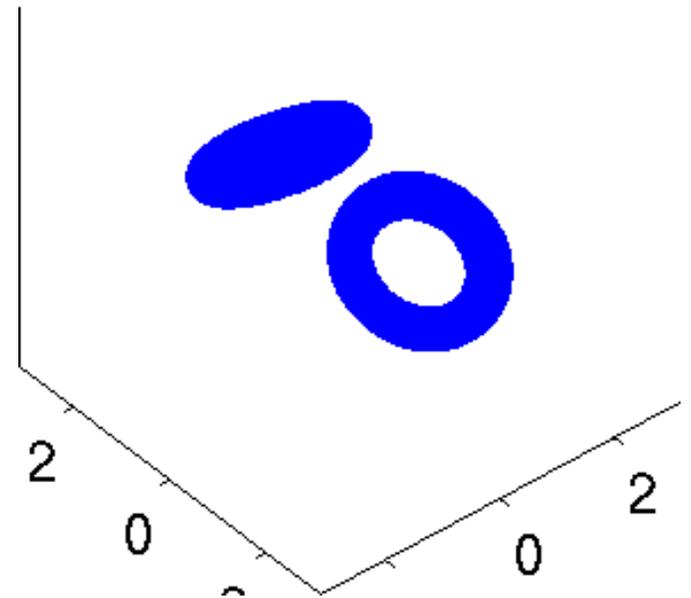
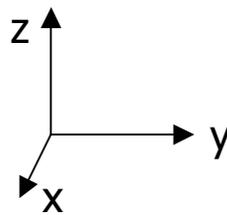
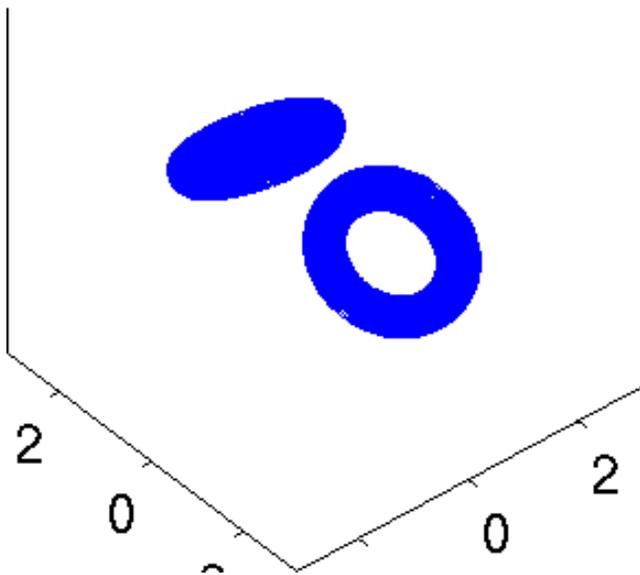
Core radius = $2 \times \text{space}$, Circulation of ring = 1

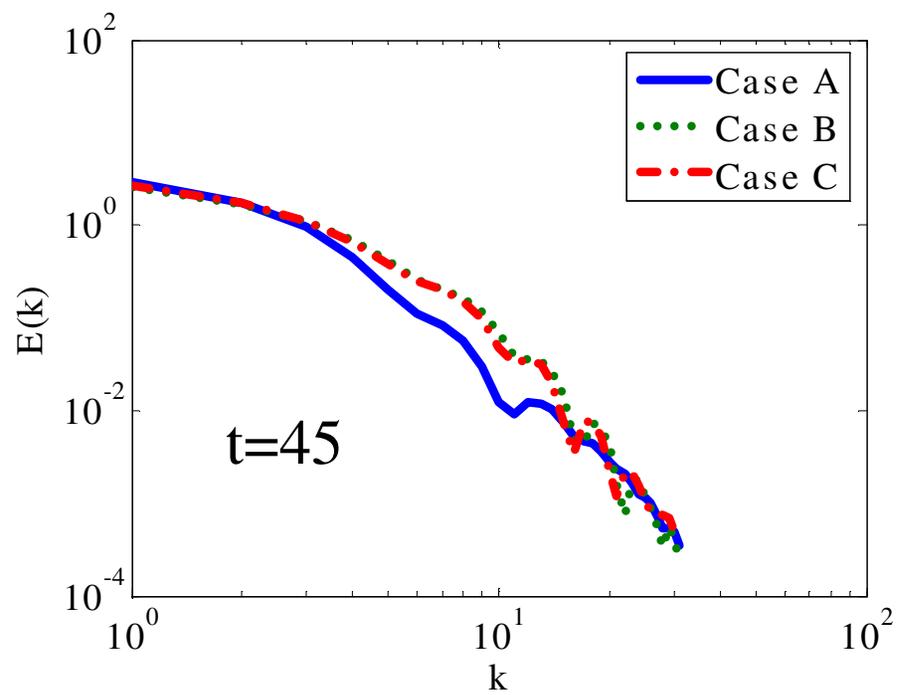
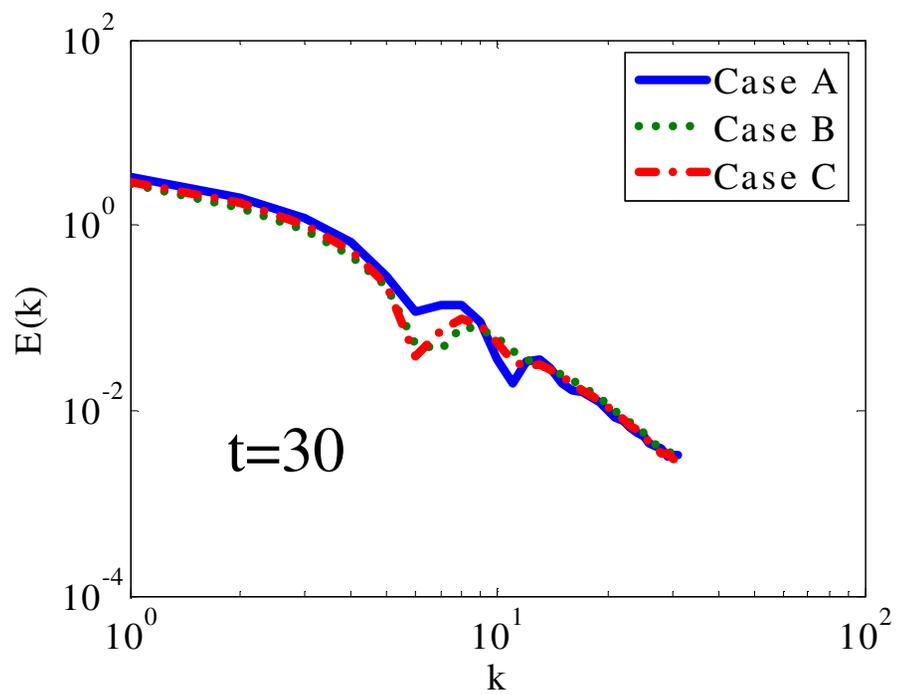
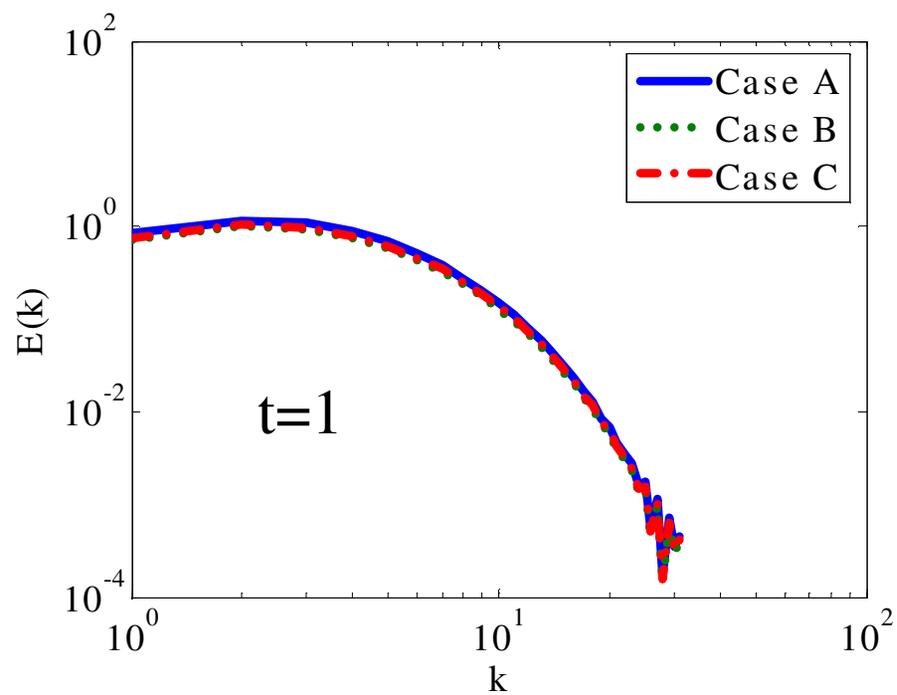
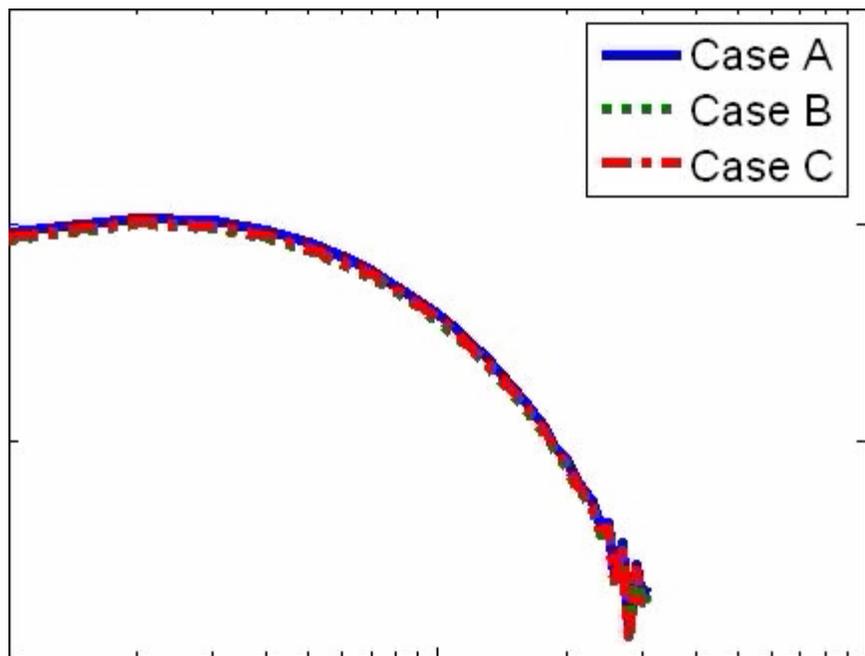
Initial distance between two rings = 3.0

Inclined angle = 30°

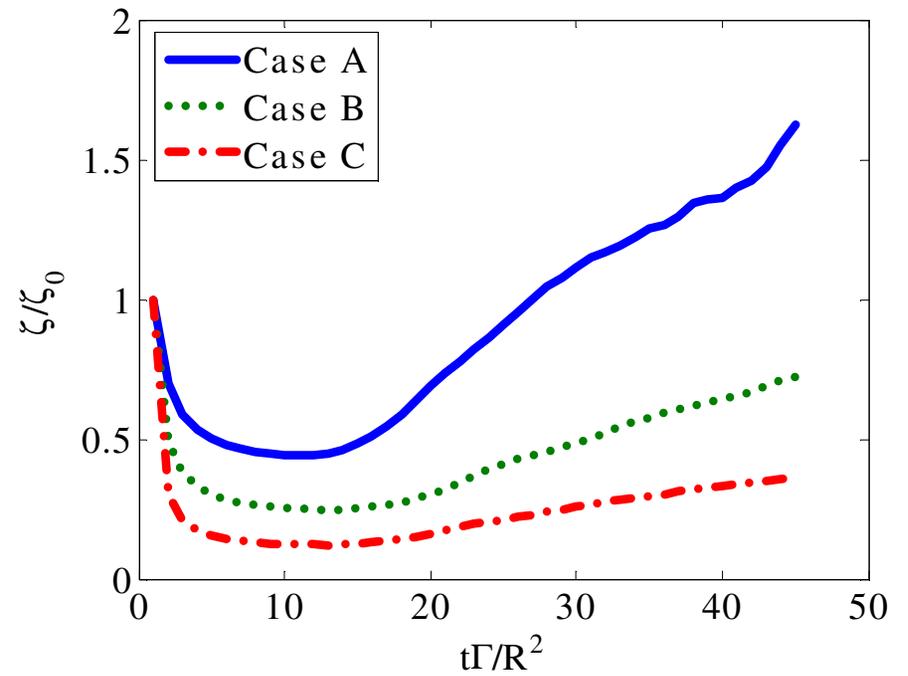
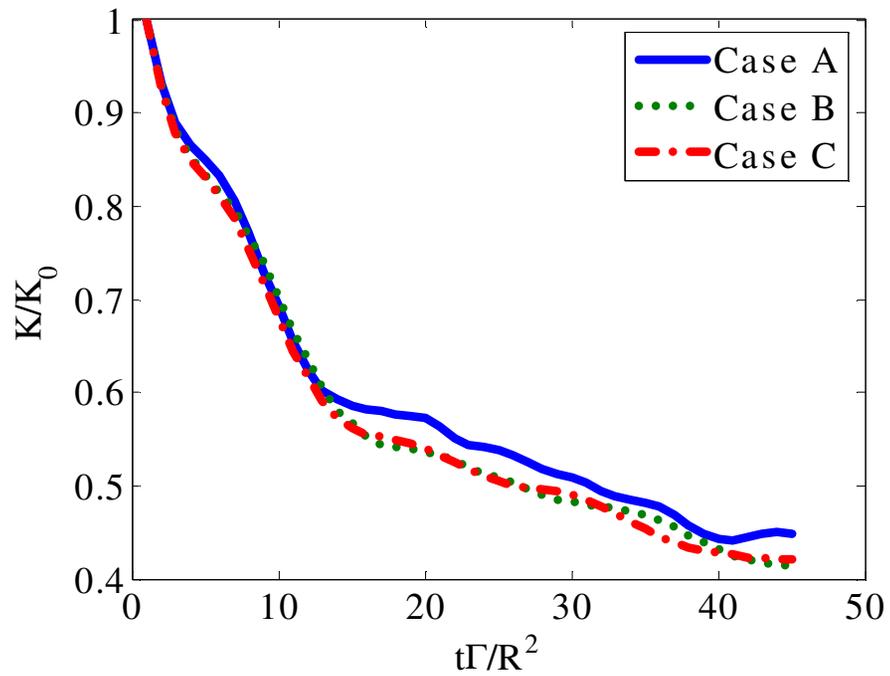
$N=10^5$

$N=10^6$





Kinetic Energy and Enstrophy



Summary

- The vortex method calculation has been accelerated dramatically with the simultaneous use of the FMM and MDGRAPE-3
- Biot-Savart calculation
 - The FMM itself accelerates the calculation 462 times, and the simultaneous use of the MDGRAPE-3 further accelerates it 4.1 times
 - The MDGRAPE-3 can accelerate the calculation 119 times, and the simultaneous use of FMM on MDGRAPE-3 is about 16 times than that of MDGRAPE-3
- Stretching term calculation
 - The calculation cost has been reduced 613 times when used FMM, and another 2.8 times faster by the simultaneous of FMM and MDGRAPE-3
 - The MDGRAPE-3 accelerates 52 times and another 33 times when combined with the FMM
- The errors involved in the use of MDGRAPE-3 are less than the errors of the FMM, and small enough to perform an accurate VM calculation
- The effect of temporal and spatial resolutions are important for accurate calculations

6

Conclusions and Outlook

Conclusion

- A fast vortex method has been developed using special-purpose computers and the simultaneous use of FMM and MDGRAPE-3
- The calculation cost has been reduced significantly by using the proposed acceleration techniques
- The reconnection of the vortex rings was clearly observed, and the discretization error became nearly negligible for the calculation using 10^7 elements
- The overall accuracy are satisfactory for VM calculations

Outlook

- The present results indicate that the calculation of further Reynolds number, the accurate vortex methods requires significantly larger N , which is possible by using the present acceleration method
- There are still some rooms to improve the acceleration rate
 - By reconstructing the subroutines which call MDGRAPE library
 - Present routines require to call MDGRAPE library in 18 times for one cross product term calculation
 - The acceleration can be improved by reducing the CALLing times
- The overall accuracy can be improved by regenerated a sophisticated function table for respective problems
- Other than the present flow, this method can be applied to calculate the homogeneous shear flow, smoothed particle hydrodynamics, dissipative particle dynamics and so on.