

To simulate turbulent thermal plasma flows for nanopowder fabrication

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Abstract: This paper discusses the theoretical models and numerical methods to simulate a turbulent thermal plasma flow transporting nanopowder growing collectively. Turbulence models should be capable of expressing not only turbulent but also laminar states because both states co-exist in and around thermal plasmas that have large variations of density as well as transport properties. Demonstrative simulations clearly present that an advanced numerical method with high-order accuracy should be used for long and robust computation capturing steep gradients of nanopowder concentration and plasma temperature and 3D dynamic motions of multi-scale vortices which are turbulent features of thermal plasma flows with low Mach numbers. High-temperature plasma regions exhibit less turbulent structures with only large vortices, whereas low-temperature regions tend to be more turbulent with numerous small vortices. In particular, a direct-current thermal plasma jet generates a double-layer structure of inner high-temperature thick vortex rings and outer low-temperature thin vortex rings near the nozzle exit. As these vortices flow downstream, they interact, deform and break up; consequently, the plasma transits to a complex thermal flow. The widely spreading distribution of the multi-scale vortices agree with an experimental observation, which is not simulated by the conventional method. Nanopowder is generated from material vapour by nucleation and condensation at the interfacial regions between plasma and cold gas. Because those regions include numerous vortices, the nanopowder is conveyed by the vortices, which results in a complex distribution of nanopowder. Simultaneously, the nanopowder diffuses and increases the size, decreasing the number by coagulation.

Keywords: Thermal plasma, Turbulent flow, Nanopowder, Numerical simulation.

1. Introduction

Thermal plasma offers a unique flow with high temperatures of not only electrons but also heavy species, high chemical reactivity, intensive light emission, and variable properties. Because it is also electrically conductive fluid, it responds to an electromagnetic field. Therefore, its plasma state can be sustained by Joule heating and controlled by Lorentz force [1-5]. Thermal plasma is anticipated as a promising device for nanopowder fabrication with remarkably high rates [6].

However, the precise process control is still difficult because the nanopowder's growth and transport must be controlled simultaneously. The process involving collective growth by nucleation, condensation, and coagulation and transport by convection, diffusion, and thermophoresis in/around a plasma is a complicated mass transfer problem in the microsecond to millisecond time scales. At present, it seems impossible to observe or measure such a spatially and temporally multi-scale process. Meanwhile, theoretical and numerical studies are potent approaches to tackle this problem, although they are arduous as well.

Turbulence of thermal plasma flows has been recognized as an important matter. Turbulence offers strong mixing effects on momentum, energy, and material molecules and particles because of multi-scale vortices. Pfender *et al.* [7] proved that a thermal plasma jet forms vortices entraining ambient cold gas at its fringe. These vortices break into smaller ones. Eventually the plasma jet transits to a

turbulent state with 3D and multi-scale vortices. For nanopowder fabrication, nanoparticles are generated not in a high-temperature plasma core but in interfacial regions between plasma and cold gas at low to intermediate temperatures [8]. Lower-temperature regions tend to be more turbulent with multi-scale vortices [9, 10]. The vortices strongly affect the transport processes of nanopowder. The fundamental mechanisms of turbulence transition of thermal plasma flows should also be clarified.

This paper discusses theoretical models and numerical methods to simulate especially a turbulent thermal plasma jet transporting nanopowder. Demonstrative simulations are performed to highlight the importance of numerical methods for long and robust computation capturing steep gradients of nanopowder concentration and plasma temperature and 3D dynamic motions of turbulent multi-scale vortices in/around a thermal plasma jet.

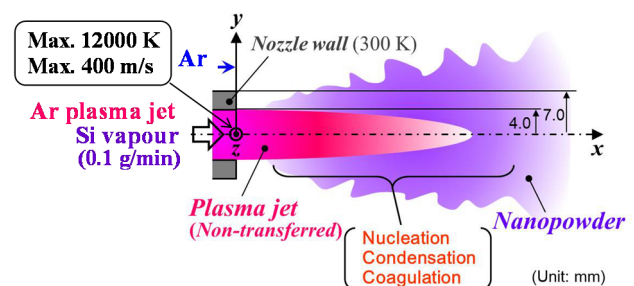


Fig. 1. A thermal plasma flow for nanopowder fabrication.

2. Model description

Thermal plasma is described by the thermofluid approximation with several typical assumptions: (i) the entire fluid region including plasma and non-ionized gas is in a local thermodynamic equilibrium state, (ii) the plasma is optically thin, and (iii) buoyancy due to density variation is negligible. The governing equations are given as the conservations of mass, momentum and energy:

$$\frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{u}) = 0 \quad (1)$$

$$\frac{\partial}{\partial t}(\rho \mathbf{u}) + \nabla \cdot (\rho \mathbf{u} \mathbf{u}) + \nabla P - \nabla \cdot \left\{ \eta \left[2\mathbf{S} - \frac{2}{3}(\nabla \cdot \mathbf{u})\mathbf{I} \right] \right\} = 0 \quad (2)$$

$$\begin{aligned} \frac{\partial}{\partial t}(\rho h) + \nabla \cdot (\rho \mathbf{u} h) - \nabla \cdot \left(\frac{\lambda}{C_p} \nabla h \right) \\ - \frac{\partial P}{\partial t} - \mathbf{u} \cdot \nabla P + Q_{rad} - Q_{con} - \Phi = 0 \end{aligned} \quad (3)$$

where ρ stands for the density of fluid, t represents the time, $\mathbf{u}$ is the velocity vector, P denotes the pressure, η signifies the viscosity, $\mathbf{I}$ is the unit matrix, h is the enthalpy, λ is the thermal conductivity, C_p is the specific heat at constant pressure, Q_{rad} represents the radiation loss, Q_{con} denotes the heat generation attributable to condensation, and Φ denotes the viscous dissipation. Furthermore, $\mathbf{S}$ is the velocity strain tensor defined as

$$\mathbf{S} = \frac{1}{2} \left[(\nabla \mathbf{u}) + (\nabla \mathbf{u})^T \right] \quad (4)$$

where superscript tr denotes the transposition.

In engineering time- and spatial-scales, the aerosol dynamics approach effectively describes the growth and transport processes of nanopowder produced by thermal plasma with several assumptions: (i) nanopowder consists of spherical nanoparticles, (ii) electric charge effects are neglected, (iii) nanoparticle temperature is identical to the fluid temperature, and (iv) material vapour is treated as an ideal gas. Extending the previous model [11] for nanopowder growth through homogeneous nucleation, heterogeneous condensation and coagulation between nanoparticles, the governing equations that also express the nanopowder's transport by convection, diffusion and thermophoresis are written as [5, 8]

$$\rho \frac{\partial}{\partial t} \left(\frac{n_p}{\rho} \right) + \rho \mathbf{u} \cdot \nabla \left(\frac{n_p}{\rho} \right) - \nabla \cdot \left[\rho D_p \nabla \left(\frac{n_p}{\rho} \right) \right] \quad (5)$$

$$- J + 2\sqrt{2}\beta_0 n_p^{11/6} f^{1/6} + \nabla \cdot \left(K_{th} \eta \frac{n_p}{\rho} \nabla \ln T \right) = 0$$

$$\rho \frac{\partial}{\partial t} \left(\frac{f}{\rho} \right) + \rho \mathbf{u} \cdot \nabla \left(\frac{f}{\rho} \right) - \nabla \cdot \left[\rho D_p \nabla \left(\frac{f}{\rho} \right) \right] \quad (6)$$

$$- J g_c - \beta_0 (n_v - n_s) n_p^{1/3} f^{2/3} + \nabla \cdot \left(K_{th} \eta \frac{f}{\rho} \nabla \ln T \right) = 0$$

$$\begin{aligned} \rho \frac{\partial}{\partial t} \left(\frac{n_v}{\rho} \right) + \rho \mathbf{u} \cdot \nabla \left(\frac{n_v}{\rho} \right) - \nabla \cdot \left[\rho D_v \nabla \left(\frac{n_v}{\rho} \right) \right] \\ + J g_c + \beta_0 (n_v - n_s) n_p^{1/3} f^{2/3} = 0 \end{aligned} \quad (7)$$

where n is the number density, D is the diffusion coefficient, and T is the temperature. The subscripts p , v , and s denote particle, vapour, and saturated state, respectively. The variable f is defined as $f = n_p g$, where g is the average monomer number in a nanoparticle. β_0 is the parameter related to collision frequency [11]. J is the homogeneous nucleation rate and g_c is the number of monomers composing a nanoparticle in a critical state [12]. K_{th} is the thermophoresis coefficient [13].

3. To express turbulent effects

A turbulent flow is composed of multi-scale vortices. A high-temperature argon thermal plasma flow cannot have sub-millimetre vortices, while a low-temperature argon flow can have vortices larger than sub-millimetres [9, 10]. Simulation of a turbulent flow of thermal plasma interacting with non-ionized cold gas requires a theoretical model and a numerical method that can express such a field.

This study adopts a Large Eddy Simulation (LES) approach with Coherent Structure (CS) model [14, 15] to deal with laminar (non-turbulent) regions and turbulent regions in and around plasma at the same time. The detailed discussion for the mathematical expression with the transformed equations is found in Refs. [8, 10].

The widely ranging temperature causes large variations of the transport properties and density with several orders of magnitude. Meanwhile, the Mach number is on the order of $10^{-3} \sim 10^{-1}$. Therefore, to obtain the solution in a practically sufficient time-scale, a thermal plasma flow is calculated as an incompressible flow with temperature-dependent density. Therefore, the following requirements should be satisfied on numerical methods.

- (i) Small vortices are resolved with a large grid element size to reduce computational costs.
- (ii) Steep gradients of variables are also captured by the computational grid element width.

Table 1. Conventional method (Method-I in Ref. [10]).

Algorithm	SIMPLE [16]
Convection terms	Upwind differencing with 1st-order accuracy
Transient terms	Backward Euler scheme with 1st-order accuracy
Diffusion terms	Central differencing with 2nd-order accuracy
Source terms	Central differencing with 2nd-order accuracy

Table 2. Advanced method (Method-III in Ref. [10]).

Algorithm	IPISO [17]
Convection terms	Hybrid-upwind K-K differencing with mainly 3rd-order accuracy [18]
Transient terms	Adams-Bashforth-Moulton scheme with 3rd-order accuracy
Diffusion terms	Central differencing with 2nd-order accuracy
Source terms	Central differencing with 2nd-order accuracy

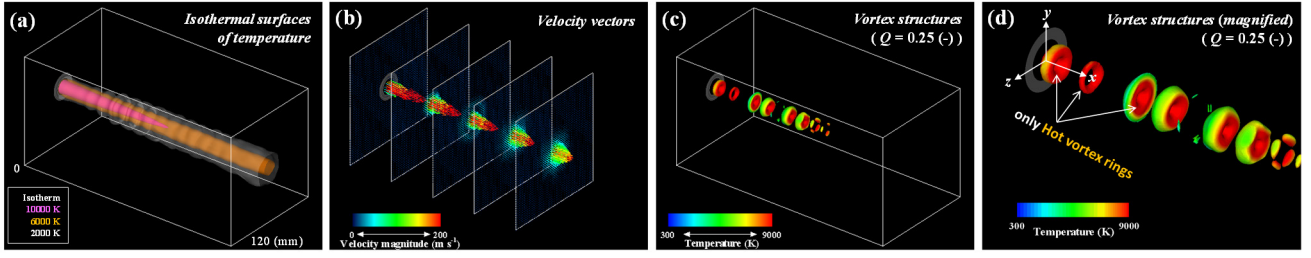


Fig. 2. Instantaneous thermal flow fields in and around a thermal plasma jet obtained by the conventional method.

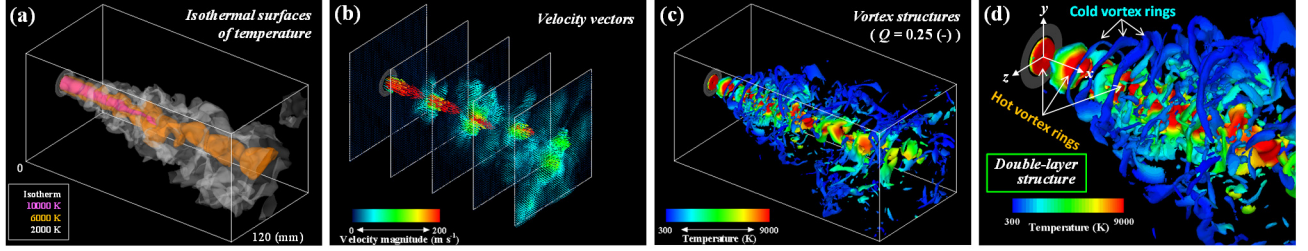


Fig. 3. Instantaneous thermal flow fields in and around a thermal plasma jet obtained by the advanced method.

- (iii) The discretization errors are small even with the large grid element width.
- (iv) The time-integration errors are small even with a large time increment.
- (v) A solution algorithm for incompressible flows is preferable even with large density variation.
- (vi) Computation is robust even with large variations of density and transport properties of plasma and extremely large values of the source terms by nanopowder formation.

Tables 1 and 2 introduce two numerical methods for the demonstrative simulations. The conventional method (Method-I [10]) has often been used for thermal plasma flow simulations because it obtains a numerical solution relatively easily. Meanwhile, the advanced method (Method-III [10]) is expected not only to express turbulent features but also to perform numerically stable computation even with a large increment of time steps. More discussion about the differencing schemes and solution algorithms is found in Refs. [8, 10].

4. Results and discussion

Figure 1 shows a schematic illustration of the computation domain. A non-transferred argon thermal plasma jet is ejected from a circular nozzle with the diameter of 8.0 mm. At the nozzle exit, the plasma jet has steady profiles of temperature with the maximum 12000 K and velocity with the maximum 400 m/s. These conditions are obtained from the approximated curves [10] with the experiment data [7]. Silicon vapour is supplied at 0.1 g/min with the plasma jet. The geometries of the domain are $(x, y, z) = (0.0 \sim 255.8 \text{ mm}, -51.2 \sim 51.1 \text{ mm}, -51.2 \sim 51.1 \text{ mm})$. Computations are performed using a Cartesian staggered grid system with the uniform spatial interval of $\Delta x = \Delta y = \Delta z = 0.1 \text{ mm}$ and the time increment of $\Delta t = 0.1 \text{ ms}$ on a supercomputer SX-ACE (NEC) at Tohoku University, Japan. The thermodynamic and transport properties of

argon thermal plasma and the material properties of silicon are obtained from Refs. [19] and [20], respectively.

Figures 2 and 3 show the snapshots of the instantaneous temperature and velocity fields and vortex structures obtained by the conventional and advanced methods, respectively. It is noted that the region near the jet nozzle $(x, y, z) = (0.0 \sim 125.0 \text{ mm}, -31.5 \sim 31.5 \text{ mm}, -31.5 \sim 31.5 \text{ mm})$ is presented here. It is apparent that the advanced method simulates a complex flow field with widely spreading multi-scale vortices. To visualize the vortices, the coherent structures of vortices were extracted from the velocity data by "*Q-criterion*" [21]. Here, the vortices are presented by the isosurfaces of the second invariant of the velocity gradient tensor of the value 0.25 which was normalized by the mean velocity of 160 m/s and the diameter of 8.0 mm at the nozzle exit. Many vortices are generated even far from the plasma jet cores. These visualized vortex structures are similar to the Schlieren photograph by Pfender *et al.* [7].

Large vortices have higher temperatures whereas small vortices have lower temperatures as predicted on the basis of Kolmogorov theory [9, 10]. The magnified snapshot Fig. 3(d) clearly shows a double-layer structure of high-temperature thicker vortex rings surrounded by low-temperature thinner vortex rings in the upstream region. As they flow downstream, the high-temperature thicker vortex rings deform largely whereas the low-temperature thinner vortex rings break up into smaller vortices.

Compared with the advanced method, the conventional method does not simulate the features of a thermal plasma jet as shown in Fig. 2. Although vortices are generated in the upstream region, they disappear in the downstream region because the upwind differencing scheme with the first-order accuracy for the convection terms inherently includes a large discretization error that unphysically flattens the spatial gradients of variables. Therefore, the

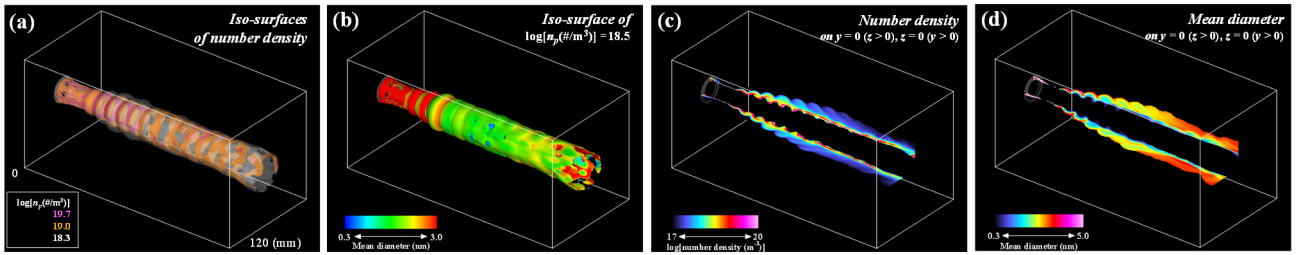


Fig. 4. Instantaneous distributions of nanopowder obtained by the conventional method.

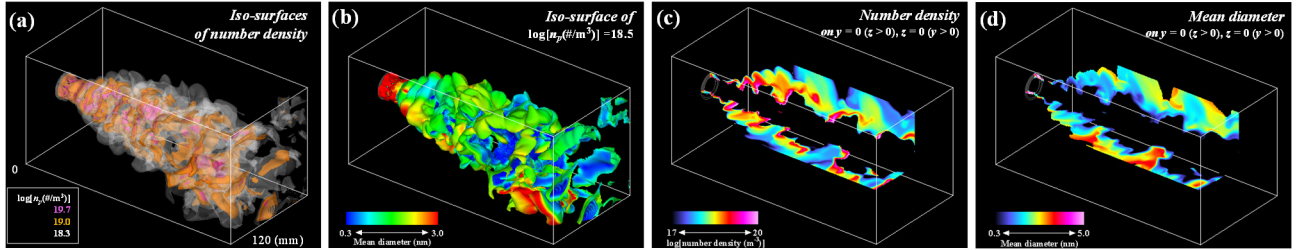


Fig. 5. Instantaneous distributions of nanopowder obtained by the advanced method.

temperature and velocity fields are calm and the vortices attenuate by the non-physical diffusional effects.

Figures 4 and 5 show the snapshots of the instantaneous number density and mean diameter distributions of silicon nanopowders obtained by the conventional and advanced methods, respectively. The silicon vapour in the plasma jet is transported by convection and diffusion. At the plasma fringe with lower temperature, the silicon vapour becomes supersaturated and changes its phase into nanopowder by homogeneous nucleation and heterogeneous condensation. Because nanopowder collectively grows up and diffuses outside the plasma region, nanopowder distributes widely in the field. The larger size regions coincide with smaller number density regions. This result indicates that simultaneous coagulation decreasing particle number plays an important role for nanopowder growth as well.

It is also obvious that the conventional method and the advanced method obtain very different results although the governing equations, the boundary conditions, and the given parameters are the same. As discussed above, the conventional method blurs gradients of variables. As a result, the vapour and the nanopowder experience unphysically large diffusions. Note, those methods obtain the very different temperature and velocity fields, which also cause the differences of the vapour and nanopowder distributions.

Using the advanced method, a numerical study has also presented that the turbulence of thermal plasma and the growth of nanopowder can be controlled by an external magnetic field [5].

5. Acknowledgements

This work was partially supported by JSPS KAKENHI (Grant No. 16K13737). The numerical results were obtained using supercomputing resources at Cyberscience Center, Tohoku University.

6. References

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