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TOWARD THE LARGE-EDDY SIMULATION OF COMPRESSIBLE TURBULENT FLOWS[†]

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ABSTRACT

New subgrid-scale models for the large-eddy simulation of compressible turbulent flows are developed and tested based on the Favre-filtered equations of motion for an ideal gas. A compressible generalization of the linear combination of the Smagorinsky model and scale-similarity model, in terms of Favre-filtered fields, is obtained for the subgrid-scale stress tensor. An analogous thermal linear combination model is also developed for the subgrid-scale models are obtained by correlating with the results of direct numerical simulations of compressible isotropic turbulence performed on a 96^3 grid using Fourier collocation methods. Extensive comparisons between the direct and modeled subgrid-scale fields are provided in order to validate the models. A large-eddy simulation of the decay of compressible isotropic turbulence — conducted on a coarse 32^3 grid — is shown to yield results that are in excellent agreement with the fine grid direct simulation. Future applications of these compressible subgrid-scale models to the large-eddy simulation of more complex supersonic flows are discussed briefly.

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1. Introduction

The direct numerical simulation of turbulent flows at the high Reynolds numbers encountered in problems of technological importance is all but impossible as a result of the wide range of scales that are present. Consequently, the solutions to such problems must invariably be based on some form of turbulence modeling. Traditional turbulence models based on Reynolds averages have had only limited success since the large scales of the turbulence – which contain most of the energy – are highly dependent on the geometry of the flow being considered. Experience has indicated that such models usually break down when a variety of turbulent flows are considered (Lumley 1983). The small scales are more universal in character, and serve mainly as a source for dissipation. Hence, it can be argued that a better understanding of turbulent flows could be achieved if just the small scales are modeled while the large scales are calculated (Deardorff 1970). This is the fundamental idea behind large-eddy simulations.

During the past decade, considerable progress has been made in the large-eddy simulation of incompressible turbulent flows. This effort has shed new light on the physics of turbulence. The earliest work relied heavily on the use of the Reynolds averaging assumption to eliminate the Leonard and cross stresses while the Reynolds stresses were computed using the Smagorinsky model (Deardorff 1970, Leonard 1974, Reynolds 1976). More recent large-eddy simulations have been based on the direct calculation of the Leonard stresses with models provided for the cross and Reynolds subgrid-scale stresses in order to enhance the numerical accuracy (see Biringen and Reynolds 1981, Bardina Ferziger and Reynolds 1983). However, among these newer models, only the Bardina, Ferziger and Reynolds (1983) model, with a Bardina constant of 1.0, satisfies the important physical constraint of Galilean invariance (Speziale 1985). The underlying physical concepts, fundamental numerical algorithms, and comprehensive historical data behind the recent field of large-eddy simulation have been presented in articles by Schumann (1975), Voke and Collins (1983) and Rogallo and Moin (1984). More recently, work on the subgrid-scale modeling of transition to turbulence of initially laminar incompressible flows has begun (Piomelli, Zang, Speziale and Hussaini 1990). Several large-eddy simulations have been performed and initial results are promising.

Despite the intensive research effort that has been devoted to the large-eddy simulation of incompressible flows as outlined above, it appears that no large-eddy simulation of a compressible turbulent flow has yet been attempted. Of course, such work could have important technological applications in the analysis of turbulent supersonic flows, where shock waves are generated, and in turbulent flows within combustion chambers. The prerequisite for carrying out such computations is the development of suitable subgrid-scale models for compressible turbulent flows. With the exception of the recent work of Yoshizawa (1986) and Speziale et al. (1988), few, if any, studies along these lines appear to have been published. The subgrid-scale models of Yoshizawa are only suitable for slightly compressible turbulent flows since they made use of an asymptotic expansion about an incompressible state. Recently however, Dahlburg, Zang and Dahlburg (1990) have performed an extensive parameter study using the model developed by Speziale et al. (1988). It was shown by Speziale et al. (1988) that, for the purposes of accuracy, the Leonard and cross stresses must be accounted for. Furthermore, the modeling of the isotropic part of the Reynolds subgrid-scale stress tensor was shown to be questionable – an issue that was left for future research.

In this paper, complete subgrid-scale models are developed for the closure of the Favrefiltered Navier-Stokes and energy equations. The compressible subgrid-scale stress model that is obtained in Section 2 reduces to the linear combination model of Bardina, Ferziger and Reynolds (1983) in the incompressible limit. Likewise, the subgrid-scale heat flux model that is obtained herein consists of an analogous linear combination of scale similarity and gradient transport terms. The dimensionless constant which appears in the subgrid-scale stress model is arrived at through correlation analysis of data generated from direct numerical simulations of compressible isotropic turbulence. A more detailed comparison of computed and modeled subgrid-scale fields is presented along with the results of a large-eddy simulation of compressible isotropic turbulence.

2. Subgrid-Scale Models for Compressible Turbulence

The compressible turbulent flow of an ideal gas is considered. Such flows are governed by the continuity, momentum and energy equations which – neglecting body forces – are given by (cf. Batchelor 1967)

$$\frac{\partial \rho}{\partial t} + \frac{\partial (\rho v_k)}{\partial x_k} = 0 \tag{1}$$

$$\frac{\partial(\rho v_k)}{\partial t} + \frac{\partial(\rho v_k v_l)}{\partial x_l} = -\frac{\partial p}{\partial x_k} + \frac{\partial \sigma_{kl}}{\partial x_l}$$
(2)

$$\frac{\partial(\rho h)}{\partial t} + \frac{\partial(\rho h v_k)}{\partial x_k} = \frac{\partial p}{\partial t} + v_k \frac{\partial p}{\partial x_k} + \frac{\partial}{\partial x_k} (\kappa \frac{\partial T}{\partial x_k}) + \Phi$$
(3)

respectively, where ρ is the mass density, **v** is the velocity vector, **p** is the thermodynamic pressure, μ is the dynamic viscosity, *h* is the enthalpy, *T* is the absolute temperature, and κ is the thermal conductivity. The viscous stress σ_{kl} and the viscous dissipation Φ are defined by

$$\sigma_{kl} = -\frac{2}{3}\mu \frac{\partial v_j}{\partial x_j} \delta_{kl} + \mu \left(\frac{\partial v_k}{\partial x_l} + \frac{\partial v_l}{\partial x_k}\right)$$
(4)

$$\Phi = -\frac{2}{3}\mu(\frac{\partial v_j}{\partial x_i})^2 + \mu(\frac{\partial v_k}{\partial x_l} + \frac{\partial v_l}{\partial x_k})\frac{\partial v_k}{\partial x_l}$$
(5)

respectively. Herein, the Einstein summation convention applies to repeated indices. Equations (1)-(3) must be supplemented with the equations of state

$$P = \rho RT, \qquad h = C_p T \tag{6}$$

for an ideal gas where R is the ideal gas constant and C_p is the specific heat at constant pressure. Likewise, the dependence of the viscosity and thermal conductivity on the temperature must be provided (i.e., relationships of the form $\mu = \mu(T)$ and $\kappa = \kappa(T)$ are needed and these depend on the gas under consideration).

Any flow variable $\mathcal F$ can be filtered in the following manner:

$$\overline{\mathcal{F}}(\mathbf{x}) = \int_D G(\mathbf{x} - \mathbf{z}, \Delta) \mathcal{F}(\mathbf{z}) d^3 z$$
(7)

where G is a filter function, Δ is the computational mesh size, and D is the domain of the fluid. The filter function G is typically taken to be an infinitely differentiable function of bounded support in a bounded domain, or a Gaussian distribution in a periodic domain[‡]. It is normalized by requiring that

$$\int_D G(\mathbf{x} - \mathbf{z}, \Delta) d^3 z = 1.$$
(8)

It follows that in the limit as the computational mesh size goes to zero, (7) becomes a Dirac delta sequence, i.e.

$$\lim_{\Delta \to 0} \int_D G(\mathbf{x} - \mathbf{z}, \Delta) \mathcal{F}(\mathbf{z}) d^3 z = \int_D \delta(\mathbf{x} - \mathbf{z}) \mathcal{F}(\mathbf{z}) d^3 z = \mathcal{F}(\mathbf{x})$$
(9)

where $\delta(\mathbf{x}-\mathbf{z})$ is the Dirac delta function (Arfken 1970). The filter function has the property that the amplitude of the high-frequency spatial Fourier components of any flow variable \mathcal{F} are substantially reduced. Consequently, $\overline{\mathcal{F}}$ represents the large-scale part of \mathcal{F} . At this point, it should be mentioned that as a result of the defining properties of G, it follows that

$$\frac{\partial \overline{\mathcal{F}}}{\partial t} = \frac{\partial \overline{\mathcal{F}}}{\partial t}, \qquad \frac{\partial \overline{\mathcal{F}}}{\partial x_{k}} = \frac{\partial \overline{\mathcal{F}}}{\partial x_{k}}.$$
(10)

Piomelli, Ferziger and Moin (1987) discuss the relationship between the form of the filter function and that of the subgrid-scale turbulence model.

The turbulent fields are decomposed as follows, based on Favre filtering:

$$\mathcal{F} = \tilde{\mathcal{F}} + \mathcal{F}' \tag{11}$$

where the Favre filter

$$\tilde{\mathcal{F}} = \frac{\overline{\rho \mathcal{F}}}{\overline{\rho}} \tag{12}$$

is defined in an analogous manner to the Favre time average which has been of use in the more traditional studies of compressible turbulent flows (Hinze 1975). However, contrary to the more traditional Favre time averaging,

$$\tilde{\tilde{\mathcal{F}}} \neq \tilde{\mathcal{F}}$$
 (13)

[‡]A Gaussian filter is adopted for the calculations in this study

in general, and hence

$$\widetilde{\mathcal{F}}' \neq 0. \tag{14}$$

The direct filtering of the continuity equation (1) yields

$$\frac{\partial \overline{\rho}}{\partial t} + \frac{\partial (\overline{\rho} \, \tilde{v}_k)}{\partial x_k} = 0 \tag{15}$$

where we have used (10) and (12). Likewise, a direct filtering of the momentum equation yields

$$\frac{\partial(\overline{\rho}\tilde{v}_{k})}{\partial t} + \frac{\partial(\overline{\rho}\tilde{v}_{k}\tilde{v}_{l})}{\partial x_{l}} = -\frac{\partial\overline{p}}{\partial x_{k}} + \frac{\partial\overline{\sigma}_{kl}}{\partial x_{l}} + \frac{\partial\tau_{kl}}{\partial x_{l}}$$
(16)

where

$$\overline{p} = \overline{\rho}R\tilde{T} \tag{17}$$

and

$$\tau_{kl} = -\overline{\rho}(\widetilde{\tilde{v}_k \tilde{v}_l} - \tilde{v}_k \tilde{v}_l + \widetilde{v'_k \tilde{v}_l} + \widetilde{v'_l \tilde{v}_k} + \widetilde{v'_k v'_l})$$
(18)

is the subgrid-scale stress tensor. The subgrid-scale stress tensor can be decomposed as follows,

$$\boldsymbol{\tau} = \mathbf{L} + \mathbf{C} + \mathbf{R} \tag{19}$$

where

$$L_{kl} = -\overline{\rho}(\widetilde{\tilde{v}_k}\widetilde{\tilde{v}_l} - \tilde{v}_k\tilde{v}_l)$$
⁽²⁰⁾

$$C_{kl} = -\overline{\rho}(v_k' \tilde{v}_l + v_l' \tilde{v}_k)$$
(21)

$$R_{kl} = -\overline{\rho} v'_k v'_l \tag{22}$$

are respectively, the subgrid-scale Leonard, cross, and Reynolds stresses based on Favre filtering. From (20), it is clear that the Leonard stress can be calculated directly and does not need to be modeled. The cross stress is modeled with the scale similarity model

$$C_{kl} = -\overline{\rho}(\tilde{v}_k \tilde{v}_l - \tilde{\tilde{v}}_k \tilde{\tilde{v}}_l) \tag{23}$$

(with a coefficient of unity to ensure Galilean invariance of the overall model). This model is analogous to its incompressible counterpart, which has been reasonably successful in the large-eddy simulation of incompressible turbulent flows (Bardina, Ferziger and Reynolds 1983, Speziale 1985). The subgrid-scale Reynolds stress tensor is separated into deviatoric and isotropic parts, respectively, as follows:

$$\mathbf{R} = {}_{D}\mathbf{R} + {}_{I}\mathbf{R} \tag{24}$$

where

$${}_{D}R_{kl} = -\overline{\rho}(\widetilde{v'_{k}v'_{l}} - \frac{1}{3}\widetilde{v'_{i}v'_{i}}\delta_{kl}), \qquad (25)$$

and

$${}_{I}R_{kl} = -\frac{1}{3}\overline{\rho}\widetilde{v'_{i}v'_{i}}\delta_{kl}.$$
(26)

Here, the deviatoric part of the subgrid-scale Reynolds stress tensor, $_{D}\mathbf{R}$, is modeled using the compressible generalization of the Smagorinsky model that is given by

$${}_{D}R_{kl} = 2C_R\overline{\rho}\Delta^2 II_{\tilde{S}}^{1/2}(\tilde{S}_{kl} - \frac{1}{3}\tilde{S}_{mm}\delta_{kl})$$

$$\tag{27}$$

where

$$\tilde{S}_{kl} = \frac{1}{2} \left(\frac{\partial \tilde{v}_k}{\partial x_l} + \frac{\partial \tilde{v}_l}{\partial x_k} \right)$$
(28)

$$II_{\mathbf{\ddot{S}}} = \tilde{S}_{mn}\tilde{S}_{mn} \tag{29}$$

(i.e., \tilde{S} is the Favre filtered rate of strain tensor while $II_{\tilde{S}}$ is its second invariant) and C_R is the compressible Smagorinsky constant. Yoshizawa – by means of a two-scale DIA method – derived a model for the isotropic part of the subgrid-scale Reynolds stress tensor, $_IR$, given by

$${}_{I}R_{kl} = -\frac{2}{3}C_{I}\overline{\rho}\Delta^{2}II_{\tilde{S}}\delta_{kl}$$

$$\tag{30}$$

where C_I is a dimensionless constant. Equation (30) can, for the most part, be obtained by making a turbulence production equals dissipation equilibrium hypothesis (Yoshizawa 1986). However, this model was shown by Speziale et al. (1988) to correlate very poorly with the results of direct numerical simulations of compressible isotropic turbulence. Since $_I\mathbf{R}$ is extremely small compared to the thermodynamic pressure, we propose to neglect it – an assumption that will be justified later. Hence, the overall subgrid-scale stress model we propose takes the form

$$\tau_{kl} = -\overline{\rho}(\widetilde{\tilde{v}_k}\widetilde{\tilde{v}_l} - \widetilde{\tilde{v}}_k\widetilde{\tilde{v}}_l) + 2C_R\overline{\rho}\Delta^2 II_{\tilde{S}}^{1/2}(\tilde{S}_{kl} - \frac{1}{3}\widetilde{S}_{mm}\delta_{kl}).$$
(31)

In the incompressible limit, Eq. (31) reduces to the linear combination model

$$-\tau_{kl}/\rho = \overline{\overline{v}_k \overline{v}_l} - \overline{\overline{v}_k} \overline{\overline{v}_l} - 2C_R \Delta^2 I I_{\overline{S}}^{1/2} \overline{S}_{kl}$$
(32)

of Bardina, Ferziger, and Reynolds (1983) where the Bardina constant is one in order to satisfy Galilean invariance (Speziale 1985). This reduction process is a consequence of Eq. (31) and

$$\tilde{\mathbf{v}} = \overline{\mathbf{v}}, \qquad \tilde{S}_{mm} = \overline{S}_{mm} = 0$$
 (33)

when ρ becomes constant.

A direct filtering of the energy equation yields the filtered form

$$\frac{\partial (C_{p}\overline{\rho}\overline{T})}{\partial t} + \frac{\partial (C_{p}\overline{\rho}\overline{v}_{k}\overline{T})}{\partial x_{k}} = \frac{\partial \overline{p}}{\partial t} + \overline{v_{k}}\frac{\partial p}{\partial x_{k}} + \overline{\Phi} + \frac{\partial}{\partial x_{k}}(\overline{\kappa}\frac{\partial T}{\partial x_{k}}) - \frac{\partial Q_{k}}{\partial x_{k}}$$
(34)

where

$$Q_{k} = C_{p}\overline{\rho} \left(\widetilde{\tilde{v}_{k}}\widetilde{T} - \widetilde{v}_{k}\widetilde{T} + \widetilde{\tilde{v}_{k}}\widetilde{T} + \widetilde{\tilde{v}_{k}}\widetilde{T}' + \widetilde{\tilde{v}_{k}}\widetilde{T}' \right)$$
(35)

is the subgrid-scale heat flux. The subgrid-scale heat flux can be decomposed in the same fashion as the subgrid-scale stresses. This leads to

$$\mathbf{Q} = \mathbf{Q}^{(L)} + \mathbf{Q}^{(C)} + \mathbf{Q}^{(R)}$$
(36)

where

$$\mathbf{Q}^{(L)} = C_{p} \overline{\rho} (\widetilde{\tilde{v}_{k}} \widetilde{T} - \widetilde{v} \widetilde{T})$$
(37)

$$\mathbf{Q}^{(C)} = C_{\mathbf{p}} \overline{\rho} (v'_{\mathbf{k}} \tilde{T} + \tilde{v}_{\mathbf{k}} \tilde{T}')$$
(38)

$$\mathbf{Q}^{(R)} = C_p \overline{\rho} \widetilde{v_k' T'} \tag{39}$$

are the Leonard, cross, and Reynolds heat fluxes. Analogous to the modeling of the cross stress, the cross heat flux is modeled using the scale similarity format

$$Q_{\boldsymbol{k}}^{(C)} = C_{\boldsymbol{p}} \overline{\rho} (\tilde{v}_{\boldsymbol{k}} \tilde{T} - \tilde{\tilde{v}}_{\boldsymbol{k}} \ \tilde{T}).$$

$$\tag{40}$$

The Reynolds heat flux is modeled with the usual gradient transport format as follows (cf. Eidson 1985):

$$\widetilde{v_k'T'} = -\frac{C_R}{Pr_T} \Delta^2 I I_{\tilde{S}}^{1/2} \frac{\partial T}{\partial x_k}$$
(41)

where Pr_T is the turbulent Prandtl number. Of course, the Leonard heat flux can be calculated directly. Hence, the overall model for the subgrid-scale heat flux we propose is as follows:

$$Q_{k} = C_{p}\overline{\rho} \left[(\widetilde{\tilde{v}_{k}}\tilde{T} - \tilde{\tilde{v}_{k}}\tilde{\tilde{T}}) - \frac{C_{R}}{Pr_{T}} \Delta^{2} I I_{\tilde{S}}^{1/2} \frac{\partial \tilde{T}}{\partial x_{k}} \right], \qquad (42)$$

and is obtained by combining equations (37), (40), and (41).

At this point, some comments need to be made concerning the viscous terms on the right-hand side of (16) and the pressure gradient-velocity and viscous dissipation terms which appear on the right-hand side of (34). The pressure gradient-velocity correlation can be written in the alternative form

$$\overline{v_{k}}\frac{\partial p}{\partial x_{k}} = \frac{\partial(\overline{pv_{k}})}{\partial x_{k}} - \overline{p}\frac{\partial v_{k}}{\partial x_{k}}$$

$$= \frac{\partial}{\partial x_{k}}(\overline{\rho RTv_{k}}) - \overline{p}\frac{\partial v_{k}}{\partial x_{k}}$$

$$= \frac{\partial}{\partial x_{k}}(\overline{\rho}R\tilde{v}_{k}\tilde{T}) + \frac{\partial}{\partial x_{k}}\left(\frac{R}{C_{p}}Q_{k}\right) - \overline{p}\frac{\partial v_{k}}{\partial x_{k}}$$
(43)

where only the pressure dilatation term

$$\overline{p}\frac{\partial v_{k}}{\partial x_{k}} = \overline{\rho}R(T\frac{\partial v_{k}}{\partial x_{k}})$$

$$= \overline{\rho}R\tilde{T}\frac{\partial \tilde{v}_{k}}{\partial x_{k}} + \overline{\rho}RT'\frac{\partial \tilde{v}_{k}}{\partial x_{k}} + \overline{\rho}R\tilde{T}\frac{\partial v'_{k}}{\partial x_{k}} + \overline{\rho}RT'\frac{\partial v'_{k}}{\partial x_{k}}$$

$$(44)$$

is not yet closed. The temperature dilatation correlation $(T' \frac{\partial v'_k}{\partial x_k})$ is extremely difficult to model and not much success has been achieved in dealing with it in the context of Reynolds stress models. However, within the framework of subgrid-scale modeling, this term and its corresponding cross correlation have physical interpretations. They represent the contribution of the dilatation of the small scales to the internal energy variation of the fluid – an effect which is expected to be small. Hence, for this initial study, we neglect such terms. Furthermore, since the mean temperature is constant and the temperature fluctuations are small (< 10%), the viscosity and thermal conductivity are held constant. For similar reasons, we also neglect the small scale component of the viscous dissipation.

The turbulence model proposed herein is thus complete once values for the constants C_R and Pr_T are obtained. This will be accomplished using the results of direct numerical simulations of compressible isotropic turbulence.

3. Numerical Method

Our direct simulations of compressible turbulence are based on a non-dimensional form of Eqs. (1)-(3), with the time derivative in the energy equation written solely in terms of the pressure. In order to alleviate the severe stability limit imposed at very low Mach numbers by the acoustic waves, a splitting method is adopted. The first step integrates the equations

$$\frac{\partial \rho}{\partial t} = 0, \tag{45}$$

$$\frac{\partial(\rho v_k)}{\partial t} + \frac{\partial(\rho v_k v_l)}{\partial x_l} = \frac{\partial \sigma_{kl}}{\partial x_l},\tag{46}$$

$$\frac{\partial p}{\partial t} + v_k \frac{\partial p}{\partial x_k} + \gamma p \frac{\partial v_k}{\partial x_k} - c_0^2 \frac{\partial (\rho v_k)}{\partial x_k} = \frac{1}{RePrM_{\infty}^2} \frac{\partial^2 T}{\partial x_k \partial x_k} + (\gamma - 1)\Phi, \qquad (47)$$

while the second step integrates

$$\frac{\partial \rho}{\partial t} + \frac{\partial (\rho v_k)}{\partial x_k} = 0, \qquad (48)$$

$$\frac{\partial(\rho v_k)}{\partial t} + \frac{\partial p}{\partial x_k} = 0, \qquad (49)$$

$$\frac{\partial p}{\partial t} + c_0^2 \frac{\partial (\rho v_k)}{\partial x_k} = 0.$$
(50)

The constants c_0 and M_{∞} are the current root mean square (rms) value of the sound speed (c) and the reference Mach number, while $\gamma = C_p/C_v$, where C_v is the specific heat at constant volume. These equations are non-dimensionalized in terms of a length scale (L_0), a velocity scale (U_0), a pressure scale (P_0), a reference viscosity μ and a reference thermal conductivity κ . The Reynolds number is given by $Re = \rho_0 U_0 L_0 / \mu$, the Prandtl number by $Pr = C_p \mu / \kappa$, and the reference Mach number by $M_{\infty} = U_0 / (\gamma R T_0)^{1/2}$. For all calculations presented in this study, $\gamma = 1.4$, and Pr = 0.7. Initially, the density ρ_0 is uniform and equal to one. The computational domain is a cube, normalized to $[0, 2\pi]^3$. Periodic boundary conditions are imposed in all three directions.

The spatial derivatives in these equations are approximated by a Fourier collocation method (see, for example, Hussaini and Zang 1987). In each coordinate direction, N grid points are used: $x_{k_j} = 2\pi j/N$, for j = 0, 1, ..., N-1. The derivative of a function $\mathcal{F}(\mathbf{x})$ with respect to x_k is approximated by the analytic derivative of the trigonometric interpolant of $\mathcal{F}(\mathbf{x})$ in the direction x_k . Most simulations of incompressible, homogeneous turbulence have used a Fourier Galerkin method. The compressible equations, however, contain cubic rather than quadratic nonlinearities and true Galerkin methods are more expensive (compared with collocation methods) than they are for incompressible flow. The essential difference between collocation and Galerkin methods is that the former are subject to both truncation and aliasing errors, whereas the latter have only truncation errors. As discussed extensively by Canuto, et al (1987), the aliasing terms are not significant for a well-resolved flow. However, care is needed to pose a collocation method in a form which ensures numerical stability. For this reason, the second term in (46) is actually used in the equivalent form

$$\frac{1}{2} \left[\frac{\partial (\rho v_k v_l)}{\partial x_l} + \rho v_l \frac{\partial v_k}{\partial x_l} + v_k \frac{\partial (\rho v_l)}{\partial x_l} \right].$$
(51)

As noted by Feiereisen, Reynolds and Ferziger (1981), when this form is employed together with a symmetric differencing method in space (for example Fourier collocation), then in addition to mass, and momentum, energy is conserved for the ideal compressible equations (zero viscosity and thermal conductivity) in the absence of time differencing (and splitting) errors.

The second fractional step of the splitting, given by (48)-(50), contains most of the effects of the acoustic waves. This splitting is employed at each stage of a third-order Runge-Kutta method. In the simulations reported here, the second fractional step is integrated analytically. In Fourier space, (48)-(50) become

$$\frac{\partial \hat{\rho}}{\partial t} + i k_l \hat{m}_l = 0 \tag{52}$$

$$\frac{\partial \hat{m}_l}{\partial t} + ik_l \hat{p} = 0 \tag{53}$$

$$\frac{\partial \hat{p}}{\partial t} + ic_0^2 k_l \hat{m}_l = 0 \tag{54}$$

where $m_l = \rho v_l$ and Fourier transformed quantities (which depend upon the wavenumber k) are denoted by a circumflex. The exact solution of these equations is

$$\hat{\rho}^{(2)} = \hat{\rho}^{(1)} + \frac{1}{c_0^2} [\hat{A} \cos(c_0 k \Delta t_s) + \hat{B} \sin(c_0 k \Delta t_s) - \hat{A}]$$
(55)

$$\hat{m}_{l}^{(2)} = \hat{m}_{l}^{(1)} - \frac{ik_{l}}{c_{0}k} [\hat{A} \sin(c_{0}k\Delta t_{s}) - \hat{B} \cos(c_{0}k\Delta t_{s}) + \hat{B}]$$
(56)

$$\hat{p}^{(2)} = \hat{A} \cos(c_0 k \Delta t_s) + \hat{B} \sin(c_0 \Delta t_s)$$
(57)

where $k = |\mathbf{k}|$ is the magnitude of the Fourier wavenumber, $\hat{A} = \hat{p}^{(1)}$, $\hat{B} = i\frac{\infty}{k}k_l\hat{m}_l^{(1)}$. The superscript 1 denotes the result of the first fractional step of the splitting and the superscript 2 the results of the second fractional step. The effective time-step of the Runge-Kutta stage is denoted by Δt_s . The advantage of this splitting is that the principal terms responsible for the acoustic waves have been isolated. Since they are treated semi-implicitly, one expects the time-step limitation to depend upon $v + |c - c_0|$ rather than v + c. (Although there is also a viscous stability limit for the first fractional step, it is well below the advection limit in the cases of interest.) This is clearly a substantial advantage at low Mach numbers. Since the second fractional step is integrated analytically, it does not contribute to any time step limitations. If one is truly interested in all the details arising from the sound waves, or if there is a substantial coupling between the sound waves and the rest of the flow, then the time-step must be small enough to resolve the temporal evolution of these waves. But, if only the larger-scale sound waves are of interest, then this splitting method is useful.

During the acoustic fractional step, an isotropic truncation is performed: for each variable $(\rho, \rho v \text{ and } p)$, all Fourier coefficients for which

$$k_l k_l \ge (N/2)^2 \tag{58}$$

are set to zero. This reduces the numerical anisotropy produced by a cubic truncation. Moreover, it reduces the aliasing interactions in the collocation method (Canuto et al 1987, Chapters 3 and 7).

The compressible code can also be executed in a purely explicit mode. In this case no splitting is performed; Eqs. (45)-(50) are simply combined in the appropriate manner and integrated directly.

The expected stability limit of this three-dimensional Fourier collocation method for the compressible Navier-Stokes equations has the form

$$\Delta t < \alpha \left[\max_{grid} \sum_{i=1}^{3} \mathcal{U}_{i} \right]^{-1} \left(\frac{2\pi}{N} \right)$$
(59)

where for the semi-implicit version,

$$\mathcal{U}_i = |v_i| + |c - c_0| \tag{60}$$

while

$$\mathcal{U}_i = |v_i| + |c| \tag{61}$$

when the time advancement is fully explicit. For the third-order Runge-Kutta method employed here, we use $\alpha = 0.5$.

A number of simulations have also been conducted of strictly incompressible flow. These were performed with a separate code which also used a Fourier collocation method, but for the simpler, incompressible Navier-Stokes equations.

4. Comparison with Incompressible Results

The initial conditions for the numerical simulations were designed to reproduce the experimental data of Comte-Bellot and Corrsin (1971) on isotropic turbulence, hereafter referred to as CBC. These experiments were also the basis of direct simulations used by Clark, Ferziger and Reynolds (1979), Bardina, Ferziger and Reynolds (1983), and McMillan and Ferziger (1979) in their analyses of incompressible LES models. Initial conditions are chosen to match CBC measurements at a non-dimensional time of 240 (cf. table 4 in Comte-Bellot and Corrsin 1971). The computational domain is a cube of side $20/2\pi$ cm. The CBC parameters are associated with measurements taken behind a grid with a mesh spacing of one inch, and a mean fluid velocity of 393.7 in/sec. The initial time in the direct simulation corresponds to t = 0.00254 sec in the CBC experiment. The reference length (L_0) , velocity (U_0) and pressure (P_0) are respectively $20/2\pi$ cm, 1 cm/sec and 1 gr/cm sec². After generating a random, divergence-free velocity profile, the kinetic energy (in Fourier space) is scaled to match the measured CBC energy spectrum (see Appendix). Finally, the velocities are scaled so that the initial rms velocity agrees with the measured values. This adjustment is typically less than 1%, which provides one measure of the uncertainty in the fit to the experimental data. With the chosen non-dimensionalization, the Reynolds number $Re = U_0 L_0 / \nu$ is 22.74 based on a kinematic viscosity $\nu = 0.14 \ cm^2/sec$. Table 1 summarizes the parameters measured

	CBC	64 ³	96 ³	128 ³
Vrms	6.75	6.75	6.75	6.75
$\frac{1}{3}tr(\mathbf{Sk})$	-	0.0	0.0	0.0
	-	68.3	68.3	68.3
ε	462	375	432	447
λ_{11}	0.26	0.28	0.27	0.25
λ_{12}	-	0.20	0.19	0.27
λ_{13}	-	0.20	0.19	0.26
R_{λ}	38.1	43.6	41.3	37.8

Table 1: Initial conditions based on CBC experiment and Clark et al. (1979) calculation. Mach number is zero.

by CBC at t=240. The Taylor microscale length λ_{kl} is defined by

$$\lambda_{kl} = \left[\frac{\langle v_k^2 \rangle}{\left\langle \left(\frac{\partial v_k}{\partial x_l}\right)^2 \right\rangle}\right]^{1/2} \tag{62}$$

and the dissipation ϵ by

$$\epsilon = 2\mu \int S_{ij} S_{ij} d^3 x, \tag{63}$$

where S_{ij} is the rate of strain tensor

$$S_{ij} = \frac{1}{2} \left(\frac{\partial v_i}{\partial x_j} + \frac{\partial v_j}{\partial x_i} \right).$$
(64)

In Eq. (62), $\langle \cdot \rangle$ denotes a spatial average. Its exact definition is given in the Appendix. The Taylor microscale Reynolds number is

$$R_{\lambda} = \frac{v_1 \lambda_{11}}{\nu}.\tag{65}$$

The velocity derivative skewness and flatness tensors Sk and Fl are the third and fourth moments of the velocity gradient and are defined by

$$Sk_{ij} = \frac{\left\langle \left(\frac{\partial v_i}{\partial x_j}\right)^3 \right\rangle}{\left\langle \left(\frac{\partial v_i}{\partial x_j}\right)^2 \right\rangle^{3/2}}$$
(66)

$$Fl_{ij} = \frac{\left\langle \left(\frac{\partial v_i}{\partial x_j}\right)^4 \right\rangle}{\left\langle \left(\frac{\partial v_i}{\partial x_j}\right)^2 \right\rangle^2}.$$
(67)

In tables 1 and 2, only the trace of the skewness and flatness tensors are shown. The remaining columns list the parameters obtained from the initial conditions of the numerical simulations on 64^3 , 96^3 and 128^3 grids. There is a 20% discrepancy between the dissipation obtained by CBC and the dissipation computed on the coarsest grid which suggests that a 64^3 grid has marginal resolution, at best. A 12% difference between the value of R_{λ} calculated on the 64^3 grid and that obtained by CBC confirms the need for grids finer than 64^3 . On a 96^3 grid, both the dissipation and R_{λ} are in much closer agreement with CBC. Discrepancies between our results and CBC for ϵ and R_{λ} are respectively 6.5% and 7.5% on a 96^3 grid. On the finest grids on which the direct simulations were performed, the computed values of ϵ and R_{λ} , respectively, have relative errors of 3.5% and less than 1% when compared to CBC.

The numerical simulations were run from t = 240 until t = 375 (in CBC units), which corresponds to a non-dimensional time interval of 0.1145 (in our units). Table 2 furnishes a comparison of the experimentally measured parameters with those from the numerical simulation at the final time. On the coarsest grid, the total dissipation rate that was calculated is still slightly below the value measured by CBC. A 96³ grid generates values of ϵ consistent with CBC.

	CBC	64 ³	96 ³	128 ³
	5.03	5.18	5.19	5.21
$\frac{1}{3}tr(\mathbf{Sk})$	-	-0.42	-0.51	-0.52
E	38.6	40.3	40.4	40.7
ε	154.4	151.3	154.6	156.8
λ_{11}	0.34	0.33	0.34	0.33
λ_{12}	-	0.24	0.24	0.23
λ_{13}	-	0.24	0.24	0.23
R_{λ}	36.6	37.8	40.4	37.2

Table 2: Final conditions (t=0.1145) based on CBC experiment and Clark et al. (1979) calculation. Mach number is zero.

At t=0.1145, the diagonal components of Sk are -0.5 which agree well with the numerical results of Kerr (1985). Kerr studied isotropic, turbulent flow, but prevented the decay of energy by using an exterior energy source at the large length scales.

As noted in the previous section, we have chosen not to de-alias the advection terms. In reaching this decision we drew upon the extensive evidence that has accumulated on aliasing effects in the last dozen years (Canuto, et al 1987., Chapters 3, 4 and 7) and upon tests conducted with the incompressible isotropic turbulence code. In this code, de-aliasing is accomplished by applying the 2/3-rule (Canuto, et al 1987, Chapters 3 and 7) in an isotropic fashion; e.g., the de-aliased results for a 64^3 grid are obtained by running the incompressible code on a 96^3 grid and applying the truncation given by (58) with N/3 in place of N/2 on the right hand side. The results are summarized in Fig. 1. Here we present the energy spectra E(k) (defined in the Appendix) for 64^3 , 96^3 , and 128^3 grids at t = 0.0586 for both aliased and de-aliased calculations. Some adverse effects of aliasing are apparent on the 64^3 grid. For the reasons outlined here, a 96^3 grid was chosen as the standard discretization for the incompressible and for the compressible simulations.

5. Compressible Turbulence Results

5.1. Direct Simulations

Recent work on the direct numerical simulation of homogeneous compressible turbulence has indicated the crucial role played by the initial conditions. Passot and Pouquet (1986) conducted direct simulations of two-dimensional, compressible isotropic turbulence and concluded that when the initial rms density fluctuations are small, the turbulence statistics remain quasi-incompressible for turbulent Mach numbers M_t less than 0.3.

$$M_t = M_{\infty} \left\langle \frac{v^2}{T} \right\rangle^{1/2}.$$
 (68)

They also demonstrated (through the use of direct numerical simulations) that eddy shocklets result for sufficiently high initial rms density fluctuations and/or turbulent Mach numbers. A more systematic analysis and categorization of the effect of the initial conditions on compressible isotropic turbulence was achieved recently by Erlebacher et al. (1990).

They concluded that for $0 < M_t < 0.3$, p_{rms} must initially be $O(M_t)$ for the resulting turbulence statistics to become strongly compressible with an O(1) ratio of compressible to incompressible turbulent kinetic energy. (For the range of M_t considered herein, no eddy-shocklets occur.) On the other hand, if initially ρ_{rms} , $T_{rms} \ll M_t$, then the resulting turbulence statistics remain quasi-incompressible.

We first present the results of direct numerical simulations of compressible isotropic turbulence corresponding to the initial conditions of the CBC experiment but with a variety of non-zero mean Mach numbers. Since for these simulations, the initial conditions are $\rho_{rms} = 0$, $T_{rms} \ll 1$, only weakly compressible turbulence statistics are expected according to the theoretical results of Erlebacher et al. (1990). Unless specified otherwise, a subscript rms for any variable \mathcal{F} refers to the quantity $\langle (\mathcal{F} - \langle \mathcal{F} \rangle)^2 \rangle^{1/2} / \langle \mathcal{F} \rangle$.

The initial pressure distribution over the entire field is specified. The fluctuating pressure, p_f is determined from the velocity distribution by enforcing a zero initial time derivative for $\nabla \cdot \mathbf{v}$. A Poisson equation for p_f is obtained from the divergence of the momentum equation after setting the time variation of $\nabla \cdot \mathbf{v}$ to zero (Feiereisen et al. 1981). The mean pressure, p_m , is then determined so that a prescribed initial mean average Mach number, M_0 , defined to be the ratio of rms fluid velocity and rms speed of sound, is achieved. An analytic expression for p_m is given by

$$p_m = \frac{M_0^2 \int v^2 d^3 x + \int \gamma p_f \rho^{-1} d^3 x}{\gamma \int \rho^{-1} d^3 x}.$$
 (69)

The initial average Mach number is specified at the outset of the direct numerical simulations (DNS) as an initial condition. Density is initially set to unity, while the temperature, if required, is derived from the equation of state. Direct numerical simulations are performed for $M_0 = 0.0, 0.1, 0.4$ and 0.6.

The Mach 0.6 case contains localized regions of supersonic flow as evidenced by tables 3-4 and by the three-dimensional contour of Mach 1 furnished in figure 2. Nonetheless, the statistical properties of the flow remain largely unaffected by compressibility effects. This is shown in figures 3-6 which track the time histories of several statistical variables obtained from 96³ DNS. The time histories for skewness (fig. 3), λ_{11} (fig. 4), and total kinetic energy (fig. 5) at Mach numbers 0.0, 0.1, 0.4 and 0.6 are virtually superimposed on each other.

Flatness and skewness are affected the most by compressibility effects in these simulations. Figure 3 indicates that the skewness which corresponds to an isotropic turbulent state monotonically increases with Mach number. It is -0.50 at $M_0 = 0$ and has increased to -0.46 at $M_0 = 0.6$. Before the flow has reached a state of isotropic turbulence, the time evolution of skewness at all Mach numbers are indistinguishable from each other. The physical system has equilibrated after approximately one third the total computation time. While not reaching an equilibrium value, it is nonetheless worthwhile to point out that the flatness parameter decreases by 2% as the Mach number is raised from 0.0 to 0.6 as seen in figure 6.

Figure 5 illustrates the decay of turbulent kinetic energy $(\int \frac{1}{2}v_iv_i d^3x)$ as a function of time. This decay is a natural consequence of viscous damping. After a brief initial increase, R_{λ} continuously decreases in time, (fig. 7), with no sign of stabilizing. On the other hand, λ which is representative of the smaller eddies, increases in time (fig. 4). This indicates that energy in the higher wavenumbers is being depleted by the molecular viscosity.

Tables 3-5 summarize the results of direct simulations of compressible isotropic turbulence for $M_0 = 0.1, 0.4$ and 0.6, for t=0.1145 on three different grids[¶]. Incompressible results are included for comparison. On all the grids, the compressible data converges to the in-

M_0	E	ε	$ abla \cdot \mathbf{v} _{max}$	$\langle \frac{1}{3} tr(\mathbf{Sk}) \rangle$	$\langle M \rangle$	M _{max}
0.0	40.26	157.4	0.00	-0.424	0.00	0.00
0.1	40.82	158.2	0.17	-0.440	0.07	0.21
0.4	41.09	160.4	1.50	-0.428	0.28	0.84
0.6	41.32	162.3	3.30	-0.406	0.43	1.26

Table 3: Summary of direct simulations on a 64^3 grid with initial average Mach numbers of 0.0, 0.1, 0.4 and 0.6 at t=0.1145.

M_0	E	ε	$ \nabla \cdot \mathbf{v} _{max}$	$\langle \frac{1}{3} tr(\mathbf{Sk}) \rangle$	$\langle M \rangle$	M _{max}
0.0	40.35	154.3	0.00	-0.506	0.00	0.00
0.1	40.49	155.5	0.14	-0.505	0.07	0.23
0.4	40.79	157.0	1.17	-0.493	0.28	0.93
0.6	41.04	158.3	2.82	-0.477	0.42	1.39

Table 4: Summary of direct simulations on a 96^3 grid with initial average Mach numbers of 0.0, 0.1, 0.4 and 0.6 at t=0.1145.

compressible results as the Mach number is driven towards zero. As expected, the divergence of velocity no longer vanishes, and is now an increasing function of M_0 .

[¶]A DNS at $M_0 = 0.6$ was not performed on the 128³ grid.

M_0	E	ε	$ \nabla \cdot \mathbf{v} _{max}$	$\langle \frac{1}{3} tr(\mathbf{Sk}) \rangle$	$\langle M \rangle$	M _{max}
0.0	40.29	153.98	0.00	-0.521	0.00	0.00
0.1	40.39	154.84	0.14	-0.518	0.07	0.21
0.4	40.78	156.54	1.11	-0.505	0.28	0.86

Table 5: Summary of direct simulations on a 128^3 grid with initial average Mach numbers of 0.0, 0.1 and 0.4 at t=0.1145

While the dissipation is approximately the same on the two finer grids, the consistently lower values on the coarsest grid confirm the previously stated conclusion that a 64^3 grid cannot resolve all the length scales. As a function of increasing Mach number, the trace of Sk increases, the dissipation decreases, while the total kinetic energy increases very slightly.

The results in tables 3-5 are averages over several DNS runs with different initial seeds. A given seed uniquely determines the initial velocity distribution, and therefore the pressure and temperature fields. Variations of the seed are only necessary to eliminate the statistical uncertainty due to the random velocity distribution. The distribution of velocity on two different grid sizes are different even when the initial seed is the same.

Skewness is even more sensitive to the grid refinement than the dissipation as witnessed by its decrease from a value of -0.505 to one of -0.521 on 96^3 and 128^3 grids respectively. This might be a result of the greater sensitivity of the fluctuating velocity field spatial derivatives to slight inaccuracies in the flow variables.

5.2. Data Analysis

Using the data generated from the previously discussed DNS of compressible homogeneous turbulence at low Mach numbers, the proposed subgrid-scale (SGS) model is now validated. Models relate the subgrid-scale stresses – which are not available to a large-eddy simulation code – to the large scale velocities which are known. These velocities are simply the Favrefiltered velocities introduced earlier. The Favre-filtered velocities are calculated by filtering the resolved DNS velocity field with a Gaussian spatial filter of width $\Delta = \Delta_f \Delta x_f$, where Δx_f is the grid spacing on the fine grid. For convenience, Δ_c and Δ_f refer to the filter width Δ non-dimensionalized with respect to the coarse and fine grid spacing respectively.

Perturbed velocity fluctuations on the fine grid are the difference between the fully resolved velocity and the filtered ones, given by

$$\mathbf{v}' = \mathbf{v} - \tilde{\mathbf{v}}.\tag{70}$$

From v' and \tilde{v} , subgrid-scale stresses based on DNS, refered to as exact, are calculated.

These include the Leonard, cross and Reynolds subgrid-scale stresses given by Eqs. (20)-(22). However, these subgrid-scale stresses themselves do not directly affect the evolution of the system. The momentum equation is only influenced by the divergence of the subgrid-scale stresses (i.e., the vector level). Similarly, $\tilde{\mathbf{v}} \cdot (\nabla \cdot \boldsymbol{\tau})$ (the scalar level) is a better representation of the dissipation terms in the energy equation than are the stresses. Consequently, correlations are performed on the tensor, vector and scalar levels. Ideally, high correlations are desired on all levels.

The data analysis proceeds in multiple stages. First, the exact stresses calculated from the DNS are injected down to the coarse grid, along with the filtered velocities. The modeled subgrid-scale stresses are then calculated (excluding the model constants) on the coarse grid. Some variables must be filtered a second time (e.g. cross stress terms). Rather than calculate them on the fine grid (which is not available to the large-eddy simulation codes), a Gaussian filter is applied to $\tilde{\mathbf{v}}$ on the coarse grid with a filter width of Δ_c . Consistency between the coarse and fine filter widths is achieved by insuring that

$$\frac{\Delta_f}{\Delta_c} = \frac{N_f}{N_c} \tag{71}$$

where N_f and N_c are respectively the number of nodes along one direction of the fine and coarse grids. This guarantees that the filtering on the coarse and fine grids is performed over the same region in physical space. Derivative evaluations on both the coarse and the fine grid are based on Fourier collocation. Calculations by McMillan and Ferziger (1979) indicate that the model constants are sensitive to the accuracy of the derivative evaluations. A general trend that has been observed is that the Smagorinsky constant is lowered when derivative quantities are evaluated more accurately. Our constants are therefore expected to lie in the lower range of the values obtained by McMillan (1980).

Next, the model constant, C_R , is calculated. Unfortunately, the constants can be calculated by a wide variety of algorithms, each with its own merits. Moreover, for each algorithm, the constants can be evaluated from tensor, vector or scalar information. Therefore, criteria must be established to identify the best method. A key test is that L + C should be Galilean invariant. To make use of this fact, an additional constant, C_C , is introduced as an extra factor in the subgrid-scale cross stress model. A self-consistent method of calculating the constants must reproduce $C_C = 1$ to satisfy the Galilean invariance property stated above (Speziale 1985). Additional tests are performed on coarse grids with varying degrees of refinement which further decrease the number of choices. A thorough discussion of model constants is the subject of the next subsection.

Once a single or a multiple set of model constants have been determined, the model subgrid-scale stresses are calculated and correlated with the exact subgrid-scale stresses calculated from the DNS after injection onto the coarse grid. The correlations are performed for each type of subgrid-scale stress individually, and for the total stress (L+C+R). Strong differences in the correlation coefficients relating total stresses are noticed depending on whether or not the Leonard stresses are included. Finally, the correlations obtained from the proposed model are compared with the linear combination model, which has been shown

to be one of the best models available for incompressible isotropic turbulence. Correlation coefficients are calculated based on the two pairs of constants that are obtained from the above considerations. The set that is finally retained corresponds to the highest levels of correlation of the total stress on the vector and scalar levels. These matters are treated more completely in a later subsection.

To avoid a possible confusion of terminology when referring to variables being compared against each other, superscripts m and e are sometimes used. They respectively refer to modeled and exact (based on DNS) variables at the tensor, vector and scalar levels.

5.2.1. Model Constants

The proposed model given by equations (31) and (42) has two undetermined coefficients. The constant, C_R , is associated with the modeled subgrid-scale Reynolds stress, \mathbf{R} , while Pr_T is associated with the thermal heat flux.

Although the cross stress model has no constant associated with it, it is nonetheless multiplied by a constant C_C . This is done in the hope of reducing the number of schemes by which the constants can be calculated. A good model should reproduce a cross stress constant of one to guarantee Galilean invariance. Once the constants have been determined, C_C is set to one and forgotten. Because the flow is isotropic, constants are expected to be the same for the three diagonal stress components, the three off-diagonal components and the three vector components. Therefore, the values presented in the tables below are averaged over the appropriate components. In the tables, D refers to averaged diagonal components, OD to averaged off-diagonal components, V to averaged vector components and S to averaged scalar components. Similar averagings are performed for the correlation coefficients.

The two constants (C_R and C_C) are calculated using two techniques – each applied on the tensor, vector and scalar levels. One method enforces equality of the rms levels of the exact and modeled stresses. This is done for each individual subgrid-scale stress (i.e., the subgrid-scale Reynolds stress and the cross stress). Hereafter this approach is referred to as RMS. The second method utilizes multiple linear least square regression (LSQ) between the exact (18) and the modeled (31) total subgrid-scale stresses to determine the constants. Table 6 summarizes the results obtained from incompressible data. The three cases presented are identical except for the initial random number seed. The constants are independent of the detailed velocity statistics. These results are based on a vector level comparison between the modeled and exact stresses. Both RMS and LSQ produce C_C close to unity as required. Unfortunately this prevents a rational choice from being made between the two approaches. A more complete set of LSQ constants is presented in table 7. They are computed at Mach numbers of 0.0, 0.1, 0.4 and 0.6 on the tensor, vector and scalar level. Computations are performed on a coarse grid of 16^3 .

	LS	Q	RM	ſS
Seed	C_R C_C C		C_R	C_{C}
1	0.012	1.04	0.023	1.03
2	0.012	1.03	0.022	1.04
3	0.012	1.03	0.022	1.03

Table 6: Model constants calculated by LSQ and RMS between exact and modeled total stresses (L+C+R). Results are based on three identical incompressible simulations except for the random initial velocity distributions. Calculations are on the vector level on a 16^3 grid.

At first glance, C_C is near unity at both the scalar and vector levels. However whereas on the vector level, the constant remains within 4% of unity for all Mach numbers, this is not the case on the scalar level where C_C is a decreasing function of Mach number. This trend is present in the data generated from both random seeds. Although not presented here, the rms cross stress model constant is also near unity when calculated based on vector level stresses. Therefore, a preferred method for the determination of the model constants is still not possible.

5.2.2. Filter Width and Grid Coarseness

Confirmation of the numerical evidence presented by McMillan and Ferziger (1979) that $\Delta_c = 2$ is the best filter width is given in table 8 ($M_0 = 0.1$). The criterion used to determine the validity of the filter width is that C_C^{RMS} must remain close to unity on the vector level. Only when $\Delta_c = 2$ is C_C near one. Similar results hold for LSQ constants. The constants also vary with respect to the coarse grid on which the LES is to be performed. Table 9 summarizes the model LSQ and RMS constants evaluated from modeled stresses calculated on 16³ and 32³ grids on the vector level. The data ($M_0 = 0.1$) shows that C_R varies by 30% when the grid size on which the modeled stresses are calculated ranges from 16³ to 32³. On the other hand, large eddy simulations using finite-difference algorithms might be performed on grids as large as 128³. Unless a subgrid-scale model is found which produces constants independent of the coarse grid size, LES simulations will run the risk of producing the wrong results. Perhaps a more complicated dependence of the modeled subgrid-scale Reynolds stresses on Δ is required.

The best model constants will produce the highest correlations between the modeled and exact subgrid-scale stresses on all levels (tensor, vector and scalar). Based on the previous discussion, only constants calculated on a vector level are adequate because they produce a C_C of unity. Unfortunately, a clear cut choice between RMS and LSQ constants cannot be made because C_C (on the vector level) is nearly one in both cases. Rather than make an arbitrary choice, both sets of constants are considered when correlating exact against modeled stresses. For reference, the constants used henceforth are

$$C_R^{LSQ} = 0.012$$
 (72)

$$C_R^{RMS} = 0.023.$$
 (73)

			See	d 1		Seed 2		
Mo		0.0	0.1	0.4	0.6	0.0	0.6	
	D	1.32	1.32	1.32	1.31	1.32	1.31	
C_{C}	V	1.04	1.04	1.02	1.00	1.04	1.00	
	S	1.00	1.00	0.95	0.6 0.0 0.6 1.31 1.32 1.31 1.00 1.04 1.00 0.873 0.971 0.934 0.018 0.016 0.015 0.012 0.012 0.012	0.934		
	D	0.018	0.018	0.018	0.018	0.016	0.015	
C_R	V	0.012	0.012	0.012	0.012	0.012	0.012	
	S	0.015	0.015	0.015	0.014	0.015	0.015	

Table 7: LSQ model constants. Filter widths are $\Delta_f = 12$ and $\Delta_c = 2$.

		LSQ			RMS		
Δ_f	6	12	24	6	12 24		
Δ_{c}	1 2		2 4 1		2	4	
C_R	0.007	0.012	0.020	0.019	0.023	0.034	
C_{C}	0.31	1.03	1.33	0.82	1.03	1.13	

Table 8: LSQ and RMS model coefficients between exact and modeled stresses on the vector level at $M_0 = 0.1$. Results are obtained with fine filter widths of 6, 12 and 24 while maintaining the proper ratio of 6 between fine and coarse widths. The coarse grid is 16^3 .

5.2.3. Correlations

Correlation coefficients have long been a preferred diagnostic tool for estimating the *reliability* of the modeled stresses. However, the subgrid-scale stresses have been separated into various components (L, C, R), each modeled separately and although the correlation of any of these components against their models might be excellent, it is still possible for the correlation of the exact total stress against the modeled total stress to be less impressive. Such is the case, for instance, when two stress components have opposite signs and partially cancel each

	LSC	5	RMS		
Grid	C_R	C_C	C_R C_C		
16 ³	0.0125	1.03	0.023	1.03	
32 ³	0.0094	1.03	0.013	1.02	

Table 9: LSQ and RMS model constants based on subgrid-scale stresses evaluated on 16^3 and 32^3 coarse grids.

other out. As a final note, before the specific correlation coefficients are presented, one must always be attentive to the actual relationship between the exact and the modeled variable, even when the correlation coefficient is relatively high (say, above 70%). A correlation coefficient as high as 70% may not be as good as it seems. For example, the correlation between the functions y = x and $y = \exp(-x)$ in the interval [0, 1] is approximately 70% although they are qualitatively different functions! As a consequence, correlations are only deemed good if the correlation coefficient is above, say, the 90% level.

For convenience, the compressible subgrid-scale model is restated here:

$${}_{D}R_{ij} = 2C_R \overline{\rho} \Delta^2 \left(\tilde{S}_{kl} \tilde{S}_{kl} \right)^{1/2} \left(\tilde{S}_{ij} - \frac{1}{3} \tilde{S}_{mm} \delta_{ij} \right)$$
(74)

$$_{I}R_{ij} = 0 \tag{75}$$

$$C_{ij} = -\overline{\rho}(\tilde{v}_i \tilde{v}_j - \tilde{\tilde{v}}_i \tilde{\tilde{v}}_j)$$
(76)

The correlation coefficients presented in table 10 between exact and modeled R and C, are independent of the model constants. The correlation coefficients are insensitive to the average Mach number variation. The neglect of $_{I}R_{ij}$ appears to be a good approximation; the direct simulations show it to be several orders of magnitude smaller than the thermodynamic pressure. For example, for all of the compressible isotropic turbulence simulations conducted in this study,

$$\frac{(\nabla \cdot {}_{I}\mathbf{R})_{rms}}{(\nabla p)_{rms}} < 3 \times 10^{-3}$$
(77)

and, hence, the effect on the isotropic part of the Reynolds stress tensor is dominated by the thermodynamic pressure.

In most of the literature on subgrid-scale models, the Leonard stress has been omitted from the total stress correlations on the grounds that it is calculated exactly (Bardina, Ferziger, and Reynolds 1983, McMillan 1980, McMillan and Ferziger 1979). However, it has recently been shown by Speziale (1985) that the combination L + C is Galilean invariant, while the Leonard and cross stresses, individually, are not. Therefore we feel that correlations of the total stress should include the Leonard stress. Table 11 summarizes the correlation coefficients between the total stress including and excluding the Leonard stress. Results

Mo		0.1	0.4	0.6
	D	31	31	31
R	OD	26	26	25
	V	22	22	22
	S	45	45	45
	D	89	89	89
C	OD	91	91	91
	V	80	80	80
	S	75	74	72

Table 10: Correlations between exact and modeled stresses, \mathbf{R} and \mathbf{C} , at Mach 0.1, 0.4 and 0.6.

are presented at Mach 0 and Mach 0.6. When the Leonard stresses are left out, correlation coefficients similar to those of Bardina are obtained on all levels. However, the inclusion of L decreases the correlations at the vector level by approximately 30%. Table 11 substantiates that the correlation coefficients (with and without L) are nearly independent of the initial average Mach number.

Correlation coefficients between $(C + R)^e$ and various combinations of the modeled stresses using constants calculated by LSQ and RMS are summarized in table 12. The second column indicates the model against which the total stress is being compared. Although at first glance RMS based constants perform better at the tensor level when the total modeled stress is considered, at the vector and scalar levels, the trend is reversed. Correlations at the vector and scalar level are higher by 4% using the LSQ constants.

When the constants are selected based on LSQ, table 12 indicates that the correlations on all levels are highest when all modeled components are included. However, the dynamic evolution of the large scale velocities only brings into play the stresses on the vector and scalar level. Therefore the coefficients which produce the maximum correlations of total stress on these two levels should be chosen. This leads to an optimum choice of

$$C_R = 0.012$$
 (78)

calculated by least squares fit of the total stress on the vector level.

Conversion of C_R to the standard form currently used in incompressible LES ¶ produces a Smagorinsky constant of 0.092. McMillan (1979) obtained a value of $C_S = 0.10$ when spectral collocation derivative computations were employed. This constant corresponds to a scalar level evaluation based on RMS. On the vector level, McMillan calculated a higher

[¶]In a number of reports (McMillan and Ferziger 1979, Bardina et al. 1983), the subgrid-scale Reynolds stress model is proportional to C_s^2 . In these cases, the relationship between C_s and C_R is $C_R = \sqrt{2}C_s^2$.

	$M_0 =$	0.0	$M_0 =$	0.6
	L+C+R	C+R	L+C+R	C+R
D	93	82	93	81
OD	80	85	79	84
V	46	72	46	71
S	56	73	56	74

Table 11: Comparison of correlation coefficients of the exact total stress versus its model. The modeled terms are computed on a 16^3 coarse grid. Each case is presented with and without the inclusion of the Leonard subgrid-scale stress terms (calculated exactly). Both the incompressible and the M=0.6 case are shown to illustrate the weak influence of Mach number on the correlation coefficients with $C_R = 0.012$.

			Least squares			RMS			
			$C_R = 0.0122$				$C_R = 1$	0.023	3
Exact	Model	D	D OD V S			D	OD	\overline{V}	S
C + R	С	78	82	68	62	78	82	68	62
	C + R	81	84	71	70	88	83	67	67

Table 12: Correlation of the exact total stress $(C + R)^e$ with various models. The modeled stresses are defined in equations (74)-(76).

Smagorinsky constant of 0.13. This value can be obtained from the present data by using ${}_{D}\mathbf{R}^{RMS}$ instead of ${}_{D}\mathbf{R}^{LSQ}$. However as shown above, the correlations of the total subgrid-scale stress would be lower.

Initial tests of the subgrid-scale heat flux model produced a turbulent Prandtl number in the range of 0.4 to 0.5. For this initial study, we take

$$Pr_T = 0.5 \tag{79}$$

which is a value that has been used in previously published large-eddy simulations of turbulent flows with thermal convection (cf. Eidson 1985). A more accurate calculation of Pr_T could be accomplished in a compressible flow with significant mean temperature gradients; this is beyond the scope of the present study and must await future research.

6. Large-Eddy Simulation of Compressible Isotropic Turbulence

Now, in order to demonstrate the efficacy of the subgrid-scale models derived in this paper, a large-eddy simulation of compressible isotropic turbulence is conducted. Since most highspeed compressible turbulent flows have significant regions where the turbulence statistics are quasi-incompressible, it is important that the models perform well for weakly compressible turbulence – the type of flow considered in the last section for the *a priori* tests. However, it is well known that *a priori* tests only provide a relatively weak gauge for the performance of subgrid-scale models in an actual large-eddy simulation (see Hussaini, Speziale and Zang 1990). It is therefore important to examine their performance in an actual large-eddy simulation, particularly for a case where significant compressibility effects are exhibited in the turbulence statistics.

Direct simulations of compressible isotropic turbulence were conducted corresponding to the initial conditions Re = 250, $\langle M_t^2 \rangle_0^{1/2} = 0.1$, $(\rho_{rms})_0 = 0$, $(T_{rms})_0 = 0.0626$ and two values of χ_0 : 0 and 0.2 which, respectively, correspond to initial values for $(R_\lambda)_0$ of 30.0 and 26.3. Here, $\chi = E^C/(E^I + E^C)$ where E^I and E^C are the incompressible and compressible parts of the turbulent kinetic energy, respectively. The direct numerical simulations were performed on a 96³ mesh. Characteristic energy and dissipation spectra associated with the DNS are shown in figures 8 (a)-(c) and 9 (a)-(c) for $\chi_0 = 0$ and $\chi_0 = 0.2$, respectively. Figure 8 (b) clearly shows that χ remains very small for all time, with only very slight modification of the energy spectrum in time. In contrast, figure 9 (b) shows an initial cascade of the spectrum towards the smaller scales, followed by a strong energy dissipation at the smaller scales of v_i^C . The dissipation spectra $(k^2 E(k))$ shown in figures 8 (a) and 9 (a) demonstrate that both the small and large scales are well resolved. Comparing figures 8 (c) and 9(c), the incompressible energy E^I is the same at t = 0 and t = 1.6, but is influenced by compressibility effects at the intermediate times. This influence is characterized by a slight decrease in E^I for $\chi_0 > 0$.

Integral properties of these two cases are plotted in figures 10 (a)-(e). These figures contrast the two runs corresponding to $\chi_0 = 0$ and $\chi_0 = 0.2$. The higher compressibility has a variety of effects on the flow. The total kinetic energy decays at a slightly slower rate, while both the skewness and the flatness are decreased. In other words, finite compressibility drives the flow more towards a Gaussian state. Both the integral scale L_I and the Taylor microscale λ_{11} become smaller as χ_0 increases. Lastly, the decay rate of the microscale Reynolds number is slower for finite χ_0 . A more detailed study of these effects awaits future research. It would be particularly useful to compute these statistical quantities based on the solenoidal and irrotational components of velocity.

As noted by Frisch and Orszag (1990), the three dimensional vorticity is formed of tubular-like structures. This is graphically represented in figure 11 which shows a volume rendering of ω^2 . Volumetric rendering is a visualization technique whereby rays are projected from the eye through the flow (which emits and absorbs light). The tubular structures of vorticity are contrasted with the spherical structure of $(\nabla \cdot \mathbf{v})^2$ shown in figure 12. This difference is to be expected since the dilatation satisfies an isotropic wave equation to lead-

ing order which shows no preferential direction. On the other hand, the vorticity stretching occurs along the axis of the vorticity vector, which then becomes a preferred direction.

The results for the direct simulation with the strongest compressibility effects ($\chi_0 = 0.2$) were filtered and injected onto a coarse 32^3 grid for comparison with an LES which was also performed on a 32^3 grid using the subgrid-scale models derived herein. In this manner, a direct comparison can be made between the results of the LES and the direct simulation for a flow where the turbulence statistics exhibit significant compressibility. The following turbulence statistics were compared:

- (1) the integrated average of $\omega_i \omega_i$ and $\left(\frac{\partial v_i}{\partial x_i}\right)^2$
- (2) the integrated average total and isotropic vortex stretching (denoted by $\omega_i S_{ij} \omega_j$ and $\omega^2 \nabla \cdot \mathbf{v}$, respectively),
- (3) the mean turbulence Mach number defined by $\langle M_t^2 \rangle^{1/2}$
- (4) the mean compressible, incompressible and total turbulent kinetic energy denoted by E^{C} , E^{I} and E_{TOT} , respectively,
- (5) the level of compressibility χ defined by $\chi = E^C/E_{TOT}$, and
- (6) the rms of the thermodynamic pressure, density, and temperature denoted by P_{rms} , ρ_{rms} , and T_{rms} , respectively.

These quantities represent a good choice of turbulence statistics to monitor the effects of compressibility.

In figure 13 (a)-(f), a direct comparison of these statistics for the DNS and LES is made for a filter width $\Delta_f = 2$. It is clear that the LES does an excellent job in reproducing the results of the DNS with the possible exception of χ . It should be noted that χ exhibits acoustic oscillations and hence is a difficult quantity to predict accurately; nonetheless, the LES yields results that are in good qualitative agreement with the direct simulation. The most striking result is how well the compressible turbulent kinetic energy and dilatational terms are captured.

It was found that a change in the filter width Δ_f to 1 or 3 – and an adjustment of the SGS model constants C_R and Pr_T of up to 25% – only led to a mild degradation of the results. However, the subgrid-scale models did play a crucial role in obtaining the accurate results shown in figure 13 (i.e., a 32³ direct simulation is substantially under-resolved). To illustrate this point, the results of a direct simulation on the coarse 32³ grid (i.e., an LES for $\Delta_f = 0$) is shown in figure 14 for the same test case. It is clear from this figure that the coarse grid DNS does a poor job in capturing the incompressible as well as the compressible turbulence statistics. Consequently, it is the adequate performance of the subgrid-scale

models in draining the proper amount of energy from the filtered fields (to account for the presence of unresolved scales) which leads to the excellent results obtained in figure 13. Considering the degree of compressibility exhibited by the turbulence statistics for the test case considered in this section, it would appear that the feasibility of the proposed subgrid-scale models has been established. A more extensive parametric study of this subgrid-scale model is under investigation by Dahlburg et al. (1990).

7. Conclusion

New subgrid-scale models for compressible turbulence have been developed and tested against the results of direct numerical simulations of compressible isotropic turbulence. These compressible subgrid-scale models, which were based on the Favre-filtered equations of motion for an ideal gas, contain two dimensionless constants and reduce to the linear combination model of Bardina, Ferziger and Reynolds (1983) in the incompressible, isothermal limit. The subgrid-scale stress model constant was found to assume the value of $C_R = 0.012$ which gives rise to correlations between the exact and modeled stresses that were above 70% on the tensor, vector and scalar levels – a correlation which compares favorably with those obtained in earlier work on the subgrid-scale modeling of incompressible turbulent flows. Another encouraging feature lies in the fact that these constants and their associated correlations were found to be comparatively insensitive to a mean Mach number in the range $0 < M_0 < 0.6$.

The results of a coarse grid 32^3 LES of compressible isotropic turbulence conducted with the subgrid-scale models derived in this paper were shown to be in excellent agreement with those obtained from a 96^3 direct simulation. These results are extremely encouraging since, for the case considered, on average 25% of the turbulent kinetic energy was compressible. Furthermore, the ability for the LES to accurately capture the dilatational statistics of the flow was quite surprising.

Future research will be directed on several fronts. The large-eddy simulation of a compressible turbulent flow with mean temperature gradients could lead to refinements in the subgrid-scale heat flux model. Furthermore, the large-eddy simulation of compressible, homogeneous shear flow could yield new insights into the performance of these subgrid-scale models. Near-wall modifications will also be implemented that allow for the large-eddy simulation of compressible, wall-bounded turbulent flows. While further improvements are still possible, we believe that the essential foundation for the large-eddy simulation of compressible turbulent flows has been established in this study. With future research, compressible LES could have a profound impact on the analysis of supersonic and hypersonic flows of aerodynamic importance.

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Appendix

Initial energy spectrum

Both the incompressible and compressible direct simulations impose a specified energy spectrum on the initial random velocity distribution. For comparative purposes, the data was obtained from tabular data found in Comte-Bellot and Corrsin (1971). They tabulate the function $E_{11}(k)$ which is related to the energy spectrum E(k) by

$$E(k) = \frac{1}{2}k^3 \frac{d}{dk} \left[\frac{1}{k} \frac{dE_{11}}{dk} \right]$$
(A1)

Unfortunately the data is noisy, so a least squares fit is performed on $\log E_{11}$, expressed as a fourth order polynomial in $\log k$. The final form obtained for E_{11} is

$$\log(E_{11}) = 2.64359 - 0.72602(\log k) - 0.32585(\log k)^{2} + 0.03525(\log k)^{3} - 0.02344(\log k)^{4}.$$
(A2)

Calculation of model constants

Before correlating the total exact subgrid-scale stress against its model, the model constants must be determined. There are many ways of accomplishing this among which two are retained. The total modeled stress is written as a linear combination of modeled terms $C_i \tau_i^m$:

$$\tau^m = \sum_{i=1}^n C_i \tau_i^m \tag{A3}$$

while the exact total stress is simply

$$\tau^{e} = \sum_{i=1}^{n} \tau_{i}^{e}. \tag{A4}$$

The unknown constants to be determined are the C_i . The first approach adopted is to calculate the root mean square of the pairs $C_i \tau_i^m$ and τ_i^e and equate them for each value of *i*. The constants thus take the values

$$C_i = \frac{\sigma_{\tau_i^e}}{\sigma_{\tau_i^m}}.$$
 (A5)

This method is referred to by RMS in the text.

A least square method applied to the total stress as a whole is an alternative approach. In this case, the norm

$$||\tau^{m} - \tau^{e}||^{2} = ||\sum_{i=1}^{n} (\tau_{i}^{e} - C_{i}\tau_{i}^{e})||^{2}$$
(A6)

is minimized with respect to the coefficients C_i . This gives rise to a linear system in the coefficients which can be solved by direct methods if the number of constants is not too large. For the subgrid-scale model considered in the text, n = 3.

Definitions

For reference purposes, several statistical definitions are provided here. All variables are defined on a three-dimensional grid and are subscripted by a single index i for convenience. The average of a function \mathcal{F}_i is

$$\langle \mathcal{F} \rangle = \frac{1}{N} \sum_{i=1}^{N} \mathcal{F}_{i}.$$
 (A7)

As a function of the average, the rms of ${\mathcal F}$ is

$$\mathcal{F}_{rms} = \sqrt{\langle (\mathcal{F} - \langle \mathcal{F} \rangle)^2 \rangle}.$$
 (A8)

Correlation coefficients are fundamental in evaluating subgrid-scale models. The correlation coefficient between two functions \mathcal{F} and \mathcal{G} is

$$C(\mathcal{F},\mathcal{G}) = \frac{\langle \mathcal{F}\mathcal{G} \rangle}{\mathcal{F}_{rms}\mathcal{G}_{rms}}.$$
 (A9)

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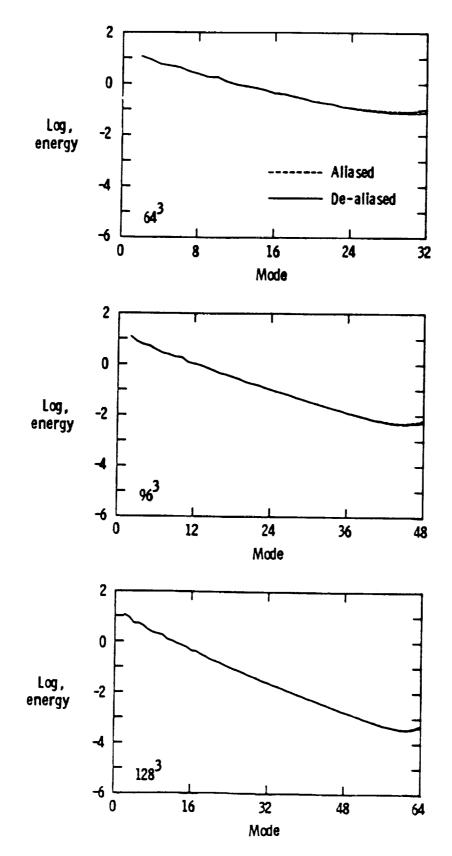


Fig. 1 Energy spectra at t = 0.0586 for both aliased and dealiased calculations for 64^3 , 96^3 , and 128^3 grids.

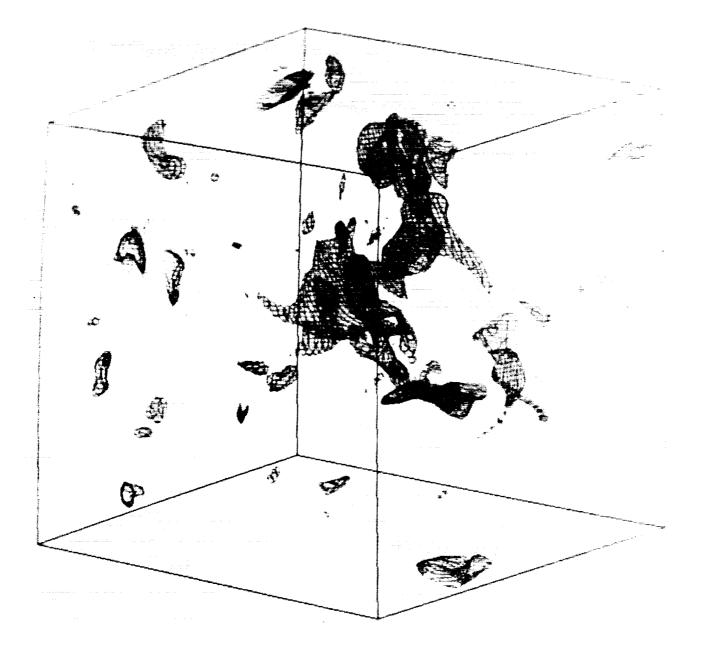


Fig. 2 Three-dimensional Mach 1.0 contours of the DNS at t = 0.0586. The direct simulation was performed on a 96³ grid.

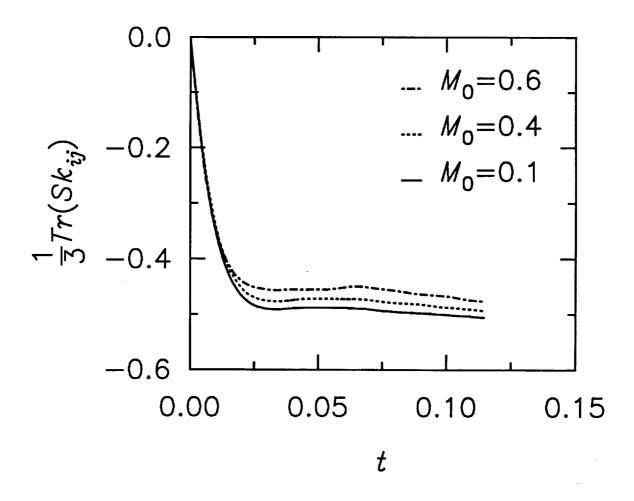


Fig. 3 Time history of one third the trace of Sk for M_0 of 0.0, 0.1, 0.4 and 0.6. Direct simulation was performed on a 96³ grid. CBC initial conditions.

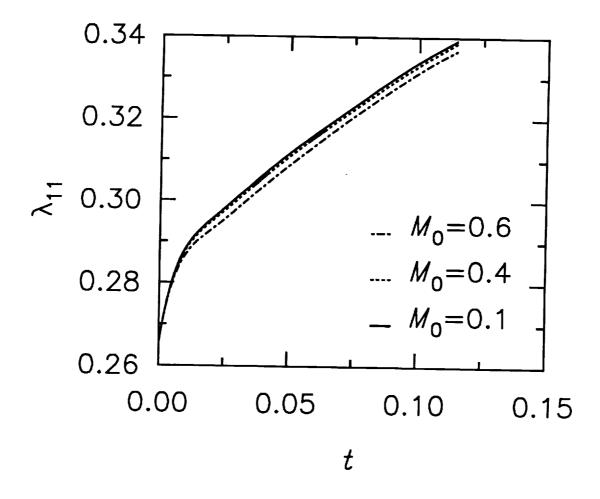


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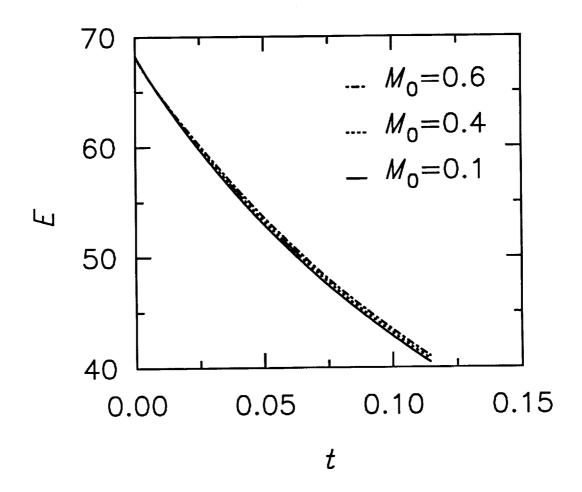


Fig. 5 Time history of kinetic energy for M_0 of 0.0, 0.1, 0.4 and 0.6. Direct simulation was performed on a 96³ grid. CBC initial conditions.

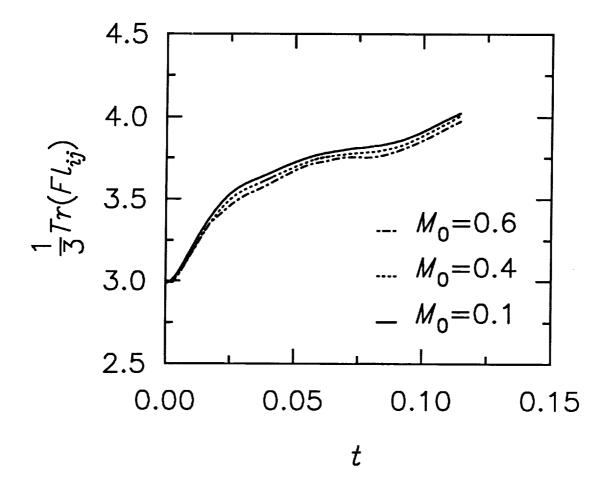


Fig. 6 Time history of one third the trace of Fl for M_0 of 0.0, 0.1, 0.4 and 0.6. Direct simulation was performed on a 96³ grid. CBC initial conditions.

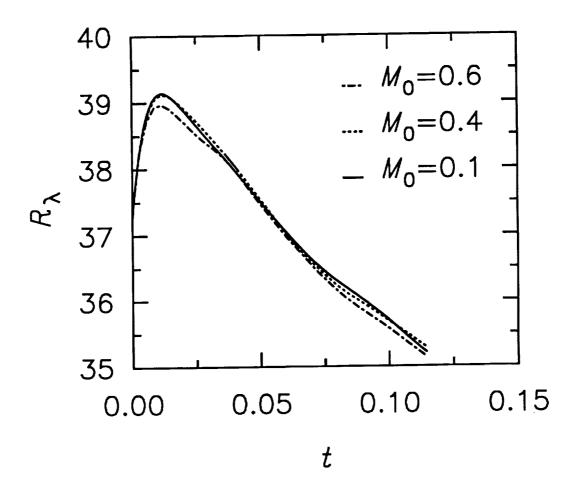
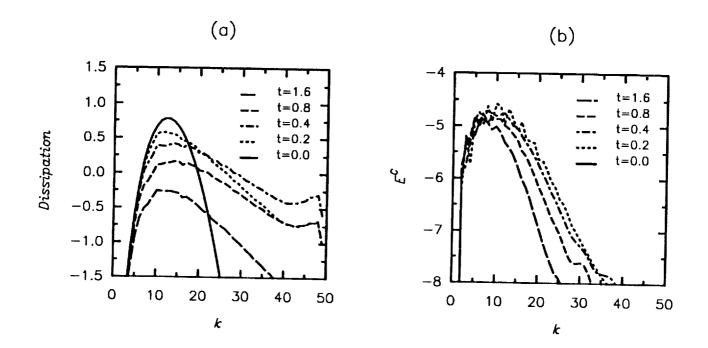


Fig. 7 Time history of R_{λ} for M_0 of 0.0, 0.1, 0.4 and 0.6. Direct simulation was performed on a 96³ grid. CBC initial conditions.



(c)

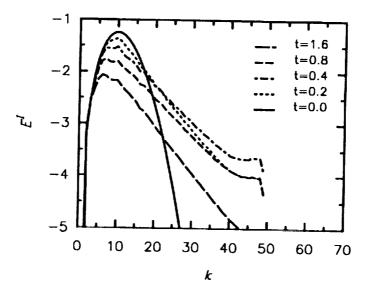
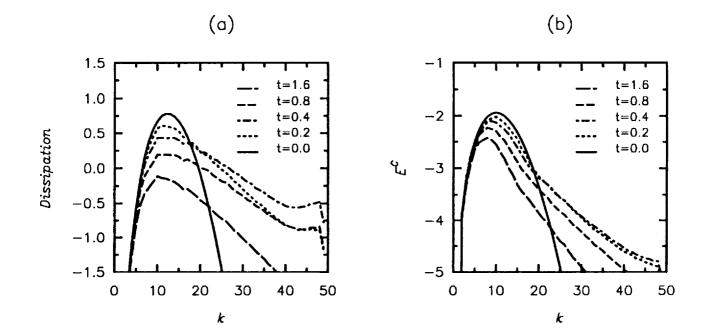


Fig. 8 Energy and dissipation spectra of 96³ DNS at t = 0, 0.2, 0.4, 0.8, 1.6. $k_0 = 10, \chi_0 = 0.0, (M_t)_0 = 0.1, (R_\lambda)_0 = 30.0, Re = 250.$ (a) Dissipation, (b) Compressible energy, (c) Solenoidal energy.



(c)

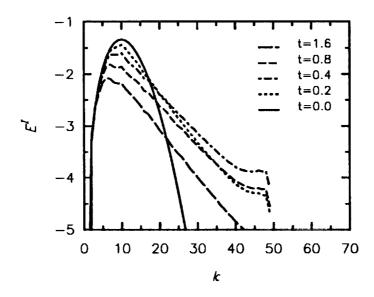


Fig. 9 Energy and dissipation spectra of 96³ DNS at t = 0, 0.2, 0.4, 0.8, 1.6. $k_0 = 10, \chi_0 = 0.2, (M_t)_0 = 0.1, (R_\lambda)_0 = 26.4, Re = 250.$ (a) Dissipation, (b) Compressible energy, (c) Solenoidal energy.

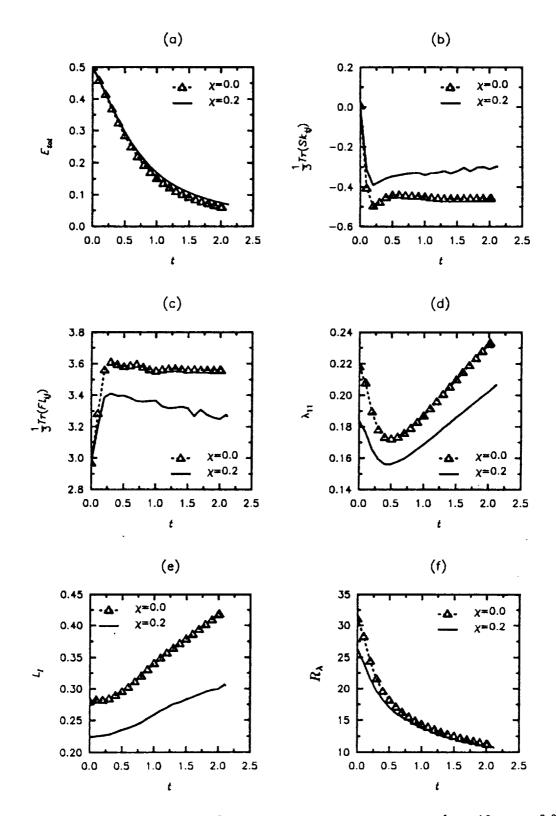


Fig. 10 Turbulence statistics for 96³ DNS at t = 0.8. Parameters are $k_0 = 10$, $\chi_0 = 0.0$ and $\chi_0 = 0.2$, $(M_t)_0 = 0.1$, $(\rho_{rms})_0 = 0$, $(T_{rms})_0 = 0.0626$, Re = 250. (a) $E_{TOT} = E^C + E^I$, (b) trace of skewness tensor, (c) trace of flatness tensor, (d) trace of Taylor length tensor, (e) integral length scale, (f) microscale Reynolds number.

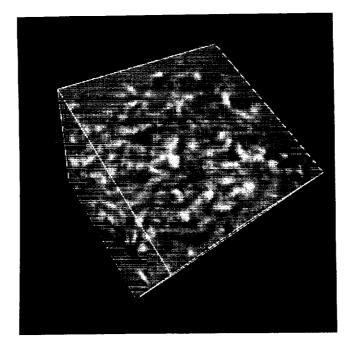


Fig. 11 Volumetric representation of ω^2 . Parameters are identical to those of Fig. 10.

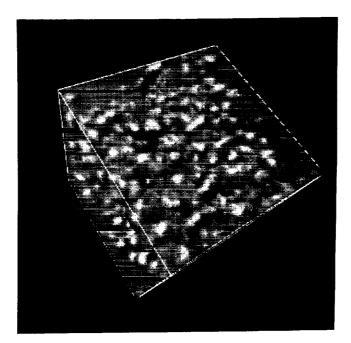


Fig. 12 Volumetric representation of $(\nabla \cdot \mathbf{v})^2$. Parameters are identical to those of Fig. 10.

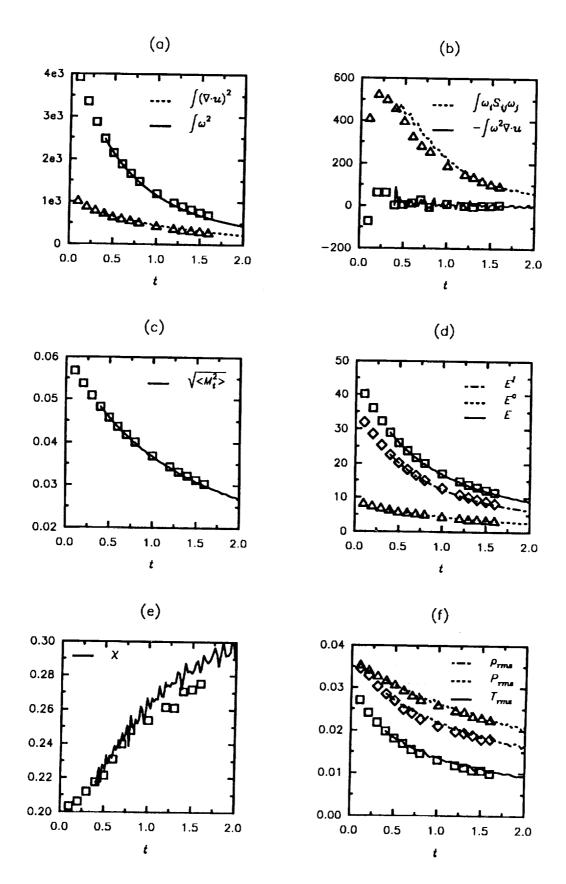
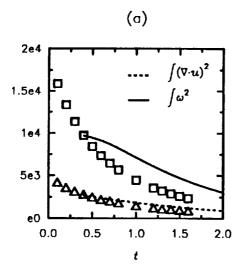
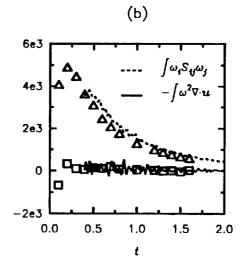
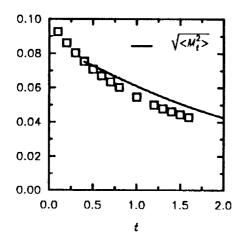


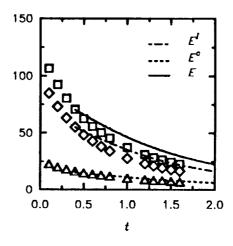
Fig.13 Comparison between LES results (---, ---) and results from a 96³ DNS $(\triangle, \Box, \diamondsuit)$ injected onto the 32³ LES grid. Filter width $\Delta_c = \Delta_f = 2$.







(c)



(d)

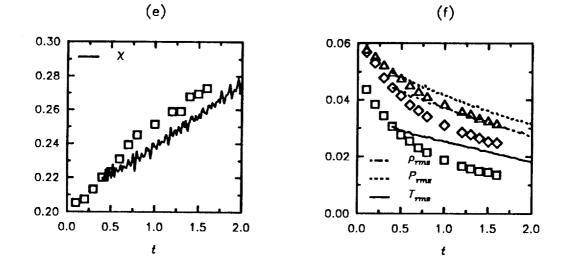


Fig.14 Comparison between LES results (---, ---) and results from a 96³ DNS (Δ, \Box, \diamond) injected onto the 32³ LES grid. Filter width $\Delta_c = \Delta_f = 0$.

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16. Abstract					
New subgrid-scale models for the large-eddy simulation of compressible turbu- lent flows are developed and tested based on the Favre-filtered equations of motion for an ideal gas. A compressible generalization of the linear combination of the Smagorinsky model and scale-similarity model, in terms of Favre-filtered fields, is obtained for the subgrid-scale stress tensor. An analogous thermal linear combina- tion model is also developed for the subgrid-scale heat flux vector. The two dimen- sionless constants associated with these subgrid-scale models are obtained by cor- relating with the results of direct numerical simulations of compressible isotropic turbulence performed on a 96 ³ grid using Fourier collocation methods. Extensive comparisons between the direct and modeled subgrid-scale fields are provided in or- der to validate the models. A large-eddy simulation of the decay of compressible isotropic turbulence conducted on a coarse 32 ³ grid is shown to yield results that are in excellent agreement with the fine grid direct simulation. Future ap- plications of these compressible subgrid-scale models to the large-eddy simulation of more complex supersonic flows are discussed briefly. 17. Key Words (Suggested by Author(s)) large-eddy simulation; compressible tur- bulence; subgrid-scale models; isotropic					
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