Drag Reduction by a Linear Viscosity Profile

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Drag reduction by polymers in turbulent flows raises an apparent contradiction: the stretching of the polymers must increase the viscosity, so why is the drag reduced? A recent theory proposed that drag reduction in agreement with experiments is consistent with the effective viscosity growing linearly with the distance from the wall. With this self-consistent solution the reduction in the Reynolds stress overwhelms the increase in viscous drag. In this Letter we show, using Direct Numerical Simulations, that a linear viscosity profile indeed reduces the drag in agreement with the theory and in close correspondence with direct simulations of the FENE-P model at the same flow conditions.

The addition of few tens of parts per million (by weight) of long-chain polymers to turbulent fluid flows in channels or pipes can bring about a reduction of the friction drag by up to 80\%\textsuperscript{1,2,3,4}. In spite of a large amount of experimental and simulational data, the fundamental mechanism has remained under debate for a long time\textsuperscript{1,2,3,4}. Since polymers tend to stretch in a turbulent flow, increasing thus the bulk viscosity, it appears contradictory that they should reduce the drag. There must exist a mechanism that compensates for the increased viscosity. Indeed, drag is caused by two reasons, one viscous, and the other inertial, related to the momentum flux from the bulk to the wall. For a fixed rate (per unit mass) of momentum generated by the pressure gradient, reducing the momentum flux can reduce the drag. In a recent theory of drag reduction in wall turbulence\textsuperscript{2} it was proposed that the polymer stretching gives rise to a self-consistent effective viscosity that increases with the distance from the wall. Such a profile reduces the Reynolds stress (i.e. the momentum flux to the wall) more than it increases the viscous drag; the result is drag reduction. The aim of this Letter is to substantiate this mechanism for drag reduction on the basis of Direct Numerical Simulations.

The onset of turbulence in channel or pipe flows increases the drag dramatically. For Newtonian flows (in which the kinematic viscosity is constant) the momentum flux is dominated by the so-called Reynolds stress, leading to a logarithmic (von-Karman) dependence of the mean velocity on the distance from the wall\textsuperscript{5}. However, with polymers, the drag reduction entails a change in the von-Karman log law such that a much higher mean velocity is achieved. In particular, for high concentrations of polymers, a regime of maximum drag reduction is attained (the “MDR asymptote”), independent of the chemical identity of the polymer\textsuperscript{2}, see Fig. 1. In a recent theoretical paper\textsuperscript{2}, the fundamental mechanism for this phenomenon was elucidated: while momentum is produced at a fixed rate by the forcing, polymer stretching results in a suppression of the momentum flux from the bulk to the wall. Accordingly the mean velocity in the channel must increase. It was shown that when the concentration of the polymers is large enough there exists a new logarithmic law for the mean velocity with a slope that fits existing numerical and experimental data. The law is universal, thus explaining the MDR asymptote.

To see how this mechanism works, consider the modified Navier-Stokes (NS) equation for the polymer solutions\textsuperscript{2,11}:

\[
\frac{\partial \mathbf{U}}{\partial t} + \mathbf{U} \cdot \nabla \mathbf{U} = -\nabla p + \nabla \cdot \mathbf{T} + \nu_0 \nabla^2 \mathbf{U},
\]

where \(\nu_0\) is the kinematic viscosity of the carrier fluid and \(\mathbf{T}\) is the extra stress tensor that is due to the polymer. Denoting the polymer end-to-end vector distance...
presented in the wall units $\mathbf{r}$, the average dimensionless extension tensor $\mathbf{R}$ is $R_{ij} \equiv \langle r_i r_j \rangle$, and the extra stress tensor is (with $\omega_{ij} \equiv \partial U_i / \partial x_j$),
\begin{equation}
\mathbf{T} = \nu_p (\mathbf{\omega} \cdot \mathbf{R} + \mathbf{R} \cdot \mathbf{\omega}^\top - \partial \mathbf{R} / \partial t - \mathbf{U} \cdot \nabla \mathbf{R}) . \tag{2}
\end{equation}
Here $\nu_p$ (proportional to the polymer concentration) is the polymeric contribution to the viscosity in the limit of zero shear. Note that expression $\mathbf{R}$ for the extra stress tensor is valid for any dumbbell model of the polymeric molecule (i.e., FENE-P, Hookean, etc.).

In $\mathbf{R}$, two simplifying approximations led to a transparent semi-quantitative theory of drag reduction. The first approximation is to ignore the fluctuations of $\mathbf{R}$ as compared to its mean and to take $\mathbf{R} \approx \langle \mathbf{R} \rangle$ in Eq. (2). Then the main contribution of the stress tensor $\mathbf{T}$ in the modified NS equation can be formally written in the form of some effective, $\mathbf{R}$-dependent tensorial viscosity. The tensorial structure of the effective viscosity describes different damping of different components of the velocity fluctuations due to the preferred orientation of polymer molecules along the mean shear. The second approximation is to ignore the preferential orientation and to replace $R_{ij} \Rightarrow R_{ij}$. These two approximations allow one to simplify the dumbbell model $\mathbf{R}$ to a modified NS equation with an effective viscosity, $\nu_0 \Rightarrow \nu$, and pressure, $p \Rightarrow P$:
\begin{equation}
\partial U_i / \partial t + U_j \nabla U_i = -\nabla \nu \cdot \mathbf{P} + \nabla [\nu (\omega_{ij} + \omega_{ji})] , \quad \nu = \nu_0 + \nu_p R , \quad P = p + \partial \nu / \partial t + \mathbf{U} \cdot \nabla \nu . \tag{3}
\end{equation}

Obviously, the polymer elongation, $\mathbf{R}$, depends on the distance from the wall, leading to the corresponding dependence of the effective viscosity. Notice that the above approximations are uncontrollable. Their verification is one of the main goals of this Letter. We demonstrate that the dynamical model Eq. (3) contains the essential properties of the full FENE-P model Eqs. (1), (2).

In Ref. $\mathbf{R}$ we modified the Reynolds closure approach to the situation in which the scalar viscosity cannot be neglected. This approach, that was justified by considering a reasonable model of the coil-stretch transition of the polymers, resulted in the analytical form of the universal MDR logarithmic profile
\begin{equation}
V^+ (y^+) = \frac{1}{\kappa} \ln \left( \frac{\nu_0}{\nu} \right) , \quad \kappa \approx 0.09 , \tag{4}
\end{equation}
presented in the wall units $\mathbf{R} \equiv L \sqrt{\nu_0 / \nu} , \quad y^+ \equiv y \mathbf{R} / L , \quad V^+ \equiv V / \sqrt{\nu_0 L} . \tag{5}
\end{equation}
Here $p^+ \equiv -\partial p / \partial x$ is fixed pressure gradients in the streamwise direction $x$, $L$ is the half-width of the channel (in the wall-normal direction $y$), and $\mathbf{R}$ is the friction Reynolds number. Equation (4) is in excellent agreement with the MDR asymptote, see Fig.1. Another conclusion of $\mathbf{R}$ is that in the MDR regime the polymer extension $\mathbf{R}(y)$ is self-adjusted in order to provide a universal linear profile of the effective viscosity
\begin{equation}
\nu_{MDR} (y) = \kappa \sqrt{\nu_0 L} y . \tag{6}
\end{equation}

Needless to say, in the viscous sublayer the viscosity is Newtonian: $\nu = \nu_0$. It should be noted that the possibility of drag reduction by increasing of the viscosity looks somewhat paradoxical: in the usual Newtonian case with constant viscosity, the drag is monotonically increasing with the viscosity. The point is that for the polymer solutions the effective viscosity is not constant anymore, it increases linearly with the distance from the wall according to Eq. (6). This point is the essential difference of the theory $\mathbf{R}$ from all previous “viscous” theories of drag reduction (see, e.g., $\mathbf{R}$). Therefore a crucial test of the theory $\mathbf{R}$ is to introduce such a linear viscosity profile to the NS Eq. (5) by hand, and see whether we observe drag reduction together with its various statistical aspects.

To this aim we simulate the effective NS Eq. (6) with proper viscosity profiles (discussed below) and show that the results are in semi-quantitative agreement with the corresponding full FENE-P DNS. All simulations were done in a domain $2\pi L \times 2L \times 1.2 \pi L$, with periodic boundary conditions in the streamwise and spanwise directions, and with no slip conditions on the walls that were separated by $2L$ in the wall-normal direction. An imposed mass flux and the same Newtonian initial conditions were used. The Reynolds number $\mathbf{R}$ (computed with the centerline velocity) was 6000 in all the runs. The grid resolution is $96 \times 129 \times 96$ for the linear viscosity profile runs and $96 \times 193 \times 96$ for the FENE-P run. The latter was done with $De = 52.7$, $\eta_p = 0.1$, $R^2_{max} = 1000$. For a definition of these parameters and details of the numerical procedure see Ref. [14].

The $y$ dependence of the scalar effective viscosity was close to being piece-wise linear along the channel height, namely $\nu = \nu_0$ for $y \leq y_1$, a linear portion with a prescribed slope for $y_1 < y \leq y_2$, and again a constant value for $y_2 < y < L$. For numerical stability this profile was smoothed out according to the differential equation
\begin{equation}
\frac{d^2 \nu}{dy^2} = \frac{C \nu_0 / \sqrt{2\pi \sigma L}}{\sqrt{2\pi \sigma L}} \left[ \exp \left( -\frac{(y - y_1)^2}{2\sigma^2} \right) - \exp \left( -\frac{(y - y_2)^2}{2\sigma^2} \right) \right] , \tag{7}
\end{equation}
integrated with initial conditions $\nu(0) = \nu_0$, and $\nu'(L) = 0$. We chose $\sigma = 0.04L$, $y_1 = L / C$, $y_2 = 3L / 4$, while $C$ is the dimensionless value of the slope. Examples of four such profiles are shown in Fig. 2. The slope of the linear part of the viscosity profile [i.e., parameter $\kappa_0$ in (6)] was varied for different runs and was smaller than the theoretically predicted value for the MDR regime. The simulations with $\nu_{MDR}(y)$ require larger $\mathbf{R}$ numbers to sustain the turbulence. Included in the figure is the flat viscosity profile of the standard Newtonian flow. Since we keep the throughput constant, the runs differ in the values of friction Reynolds numbers $\mathbf{R}^\ast \equiv \sqrt{\tau \nu_0 / \nu_0}$ where
FIG. 2: The Newtonian viscosity profile and four examples of close to linear viscosity profiles employed in the numerical simulations. Solid black line: run N, solid blue: run R, red: run S, green: T, violet: run U.

\( \tau_w \) is the average friction at the wall \( \tau_w \equiv \nu_0 dU/dy \). The decreased value of \( \text{Re}_\tau \) is a manifestation of the drag reduction measured in percentage as \( DR\% = (\tau_{w}^{N} - \tau_{w}^{E})/\tau_{w}^{N} \).

The normalized slopes, the value of \( \text{Re}_\tau \) and the percentage of drag reduction for these runs are summarized in Table I.

In Fig. 3 we show the resulting profiles of \( V_0^+(y) \) vs. \( y^+ \). The line types are chosen to correspond to those used in Fig. 2. The decrease of the drag with the increase of the slope of the viscosity profiles is obvious. Since the slopes of the viscosity profiles are smaller than needed to achieve the MDR asymptote for the corresponding \( \text{Re}_\tau \), the drag reduction occurs only in the near-wall region and the Newtonian plugs are clearly visible.

It is most interesting to compare the effect of the linear viscosity profile to the simulation of the FENE-P model in which both the throughput and \( \text{Re}_\tau \) are the same. Such a comparison was performed for the “S” viscous run, for which \( \text{Re}_\tau = 214 \) and the FENE-P run with \( \text{Re}_\tau = 212.5 \). The results are presented in the two panels of Fig. 4. In the upper panel the mean velocity profiles (symbols for FENE-P and dashed line for the linear viscosity profile) are seen to correspond very closely. The region with increased slope of the mean ve-

\[ \text{FIG. 3: The reduced mean velocity as a function of the reduced distance from the wall. The line types correspond to those used in Fig. 2.} \]

\[ \text{FIG. 4: Upper panel: the reduced mean velocity as a function of the reduced distance from the wall. Lower panel: Reynolds stresses across the channel. Continuous line: Newtonian. Dashed line: linear viscosity profile. Symbols: FENE-P.} \]

\[ \text{TABLE I: DNS parameters for effective viscosity runs.} \]

| Case | C | \( \text{Re}_\tau \) | \( DR\% \) | \( \kappa_N \) |
|------|---|----------------|---------|-----------|
| N    | 0  | 245           | -       | -         |
| R    | 8  | 227           | 13.8    | 0.035     |
| S    | 9  | 214           | 21.6    | 0.042     |
| T    | 10 | 197           | 36.9    | 0.051     |
| U    | 12 | 185           | 42.0    | 0.065     |
velocities. The increase in r.m.s. streamwise ($V^{+}_x$) and decrease in r.m.s. wall-normal ($V^{+}_y$) velocity fluctuations were observed in many experiments and numerical simulations of drag-reducing flows. In Fig. 5 we compare these quantities for FENE-P and “S” runs. Clearly, the correct trend of the r.m.s velocity fluctuations is observed, indicating that the important features of the mechanism of drag reduction are correctly captured by the model. An almost quantitative agreement is reached in the region where drag reduction actually occurs ($y/L \approx 0.1 - 0.3$).

In conclusion, we showed that the simple linear viscosity model \(\tilde{\nu}\), \(\tilde{\nu}\) faithfully demonstrates drag-reducing properties and, surprisingly enough, the amount of drag reduction increases with the increase of the slope of the viscosity profile. Even more interestingly, the behavior of objects like the Reynolds stress or the velocity fluctuations in the elastic sublayer are in close correspondence with the full FENE-P model, indicating that the mechanisms of drag reduction proposed in [7] operates similarly in both cases.

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