Fluid squeeze-out between rough surfaces: comparison of theory with experiment

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Abstract
We study the time dependence of the (average) interfacial separation between an elastic solid with a flat surface and a rigid solid with a randomly rough surface, squeezed together in a fluid. We use an analytical theory describing the fluid flow factors, based on the Persson contact mechanics theory and the Bruggeman effective medium theory, to calculate the removal of the fluid from the contacting interface of the two solids. In order to test this approach, we have performed simple squeeze-out experiments. The experimental results are compared to the theoretical predictions.

(Some figures in this article are in colour only in the electronic version)

1. Introduction

Contact mechanics between solid surfaces is the basis for understanding many tribology processes [1–7] such as friction, adhesion, wear and sealing. The two most important properties in contact mechanics are the area of real contact and the interfacial separation between the solid surfaces. For non-adhesive contact and small squeezing pressure, the average interfacial separation depends logarithmically \cite{8–10}, and the (projected) contact area linearly, on the squeezing pressure \cite{11}. Here we study how the (average) interfacial separation depends on time when an elastic solid with a flat surface is squeezed against a rigid solid with randomly rough surface in a fluid \cite{12}.

The influence of surface roughness on fluid flow at the interface between solids in stationary or sliding contact is a topic of great importance both in Nature and technology. Technological applications include leakage of seals, mixed lubrication and removal of water from the tire–road footprint. In Nature fluid removal (squeeze-out) is important for adhesion and grip between the adhesive toe pads of a tree frog or a gecko and the countersurface during rain, as well as for cell adhesion.

Almost all surfaces in Nature and most surfaces of interest in tribology have roughness on many different length scales, sometimes extending from atomic distances (∼1 nm) to the macroscopic size of the system which could be of the order of ∼1 cm. Often the roughness is fractal-like so that when a small region is magnified (in general with different magnification in the parallel and orthogonal directions) it ‘looks the same’ as the unmagnified surface.

Most objects produced in engineering have some particular macroscopic shape characterized by a radius of curvature (which may vary over the surface of the solid), e.g. the radius $R$ of a cylinder in an engine. In this case the surface may appear perfectly smooth to the naked eye but at short enough length scale, in general much smaller than $R$, the surface will exhibit strong irregularities (surface roughness). The surface roughness power spectrum $C(q)$ of such a surface exhibits a roll-off wavelength $\lambda_0 \ll R$ (related to the roll-off wavevector $q_0 = 2\pi/\lambda_0$) and therefore it appears smooth (except for the macroscopic curvature $R$) on length scales much longer than $\lambda_0$. In this case, when studying the fluid flow between two macroscopic solids, one may replace the microscopic equations of fluid dynamics with effective equations describing the average fluid flow on length scales much larger than $\lambda_0$ and which can be used to study, for example, the lubrication of the cylinder in an engine. This approach of eliminating or integrating out short length scale
degrees of freedom to obtain effective equations of motion, describing the long distance (or slow) behavior, is a very general and powerful concept often used in physics.

In the context of fluid flow at the interface between closely spaced solids with surface roughness, Patir and Cheng [13, 14] have shown how the Navier–Stokes equations of fluid dynamics can be reduced to effective equations of motion involving locally averaged fluid pressures and flow velocities. In the effective equations occur so-called flow factors, which are functions of the locally averaged interfacial surface separation. They showed how the flow factors can be determined by solving numerically the fluid flow in small rectangular units with linear sizes of the order of (or larger than) the roll-off wavelength $\lambda_0$ introduced above. In [15] one of us has developed an analytical theory for the pressure flow factors based on the Persson contact mechanics model and the Bruggeman effective medium theory to take into account the topography disorder resulting from the random roughness. We will use this theory in the calculations presented below.

This paper is organized as follows: in section 2 we briefly review the basic equations of fluid dynamics and describe some simplifications which are valid for the present case; in section 3 we describe the experimental method we have used to study the interfacial separation; and in section 4 we compare the experimental results to the theory prediction. The summary and conclusions are presented in section 5.

2. Theory

2.1. Fluid flow between solids with random surface roughness

Consider two elastic solids with randomly rough surfaces. Even if the solids are squeezed in contact, because of the surface roughness there will, in general, be non-contact regions at the interface and, if the squeezing force is not too large, there will exist non-contact channels from one side to the other side of the nominal contact region. We consider now fluid flow at the interface between the solids. We assume that the fluid is Newtonian and that the fluid velocity field $\mathbf{v}(x, t)$ satisfies the Navier–Stokes equation:

$$\frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v} = -\frac{1}{\rho} \nabla p + \nu \nabla^2 \mathbf{v}$$

where $\nu = \eta/\rho$ is the kinematic viscosity and $\rho$ is the mass density. For simplicity we will also assume an incompressible fluid so that

$$\nabla \cdot \mathbf{v} = 0.$$ 

We assume that the nonlinear term $\mathbf{v} \cdot \nabla \mathbf{v}$ can be neglected (which corresponds to small inertia and a small Reynolds number), which is usually the case in fluid flow between narrowly spaced solid walls. For simplicity we assume the lower solid to be rigid with a rough surface, while the upper solid is elastic with a flat surface. A coordinate system $xyz$ with the $xy$ plane in the surface of the lower solid and the $z$ axis pointing toward the upper solid is introduced, see figure 1. The upper solid moves with the velocity $\mathbf{v}_0$ parallel to the lower solid. Let $u(x, y, z)$ be the separation between the solid walls and assume that the slope $|\nabla u| \ll 1$. We also assume that $u/L \ll 1$, where $L$ is the linear size of the nominal contact region. In this case one expects the fluid velocity to vary slowly with the coordinates $x$ and $y$ as compared to the variation in the orthogonal direction $z$. Assuming a slow time dependence, the Navier–Stokes equations reduces to

$$\frac{\partial^2 \mathbf{v}}{\partial z^2} \approx \nabla p.$$

Here and in what follows $\mathbf{v} = (v_x, v_y)$, $\mathbf{x} = (x, y)$ and $\nabla = (\partial_x, \partial_y)$ are two-dimensional vectors. Note that $v_z \approx 0$ and that $p(x)$ is independent of $z$ to a good approximation. The solution to the equations above can be written as

$$\mathbf{v} \approx \frac{1}{2\eta} (z - u_0(x))(z - u_1(x))\nabla p + \frac{z - u_0(x)}{u_1(x) - u_0(x)} \mathbf{v}_0$$

so that $\mathbf{v} = 0$ on the solid wall $z = u_0(x)$ and $\mathbf{v} = \mathbf{v}_0$ for $z = u_1(x)$. Integrating over $z$ (from $z = u_0(x)$ to $u_1(x)$) gives the fluid flow vector

$$\mathbf{J} = -\frac{u^2(x)}{12\eta} \nabla p + \frac{1}{2} u(x) \mathbf{v}_0$$

where $u(x) = u_1(x) - u_0(x)$ is the interfacial separation at $x$. Mass conservation demands that

$$\frac{\partial u(x, t)}{\partial t} + \nabla \cdot \mathbf{J} = 0$$

where the interfacial separation $u(x, t)$ is the volume of fluid per unit area. In this last equation we have allowed for a slow time dependence of $u(x, t)$ as would be the case, for example, during fluid squeeze-out from the interfacial region between two solids.

2.2. Viscosity of confined fluids

It is well known that the viscosity of fluids at high pressures may be many orders of magnitude larger than at low pressures. Using the theory of activated processes, and assuming that a local molecular rearrangement in a fluid results in a local volume expansion, one expects an exponential dependence on the hydrostatic pressure $\eta = \eta_0 \exp(p/p_0)$, where typically (for hydrocarbons or polymer fluids) $p_0 \approx 10^5$ Pa (see, e.g., [5, 16]). Here we are interested in (wetting) fluids confined between the surfaces of elastically soft solids, e.g. rubber or
gelatin. In this case the pressure at the interface is usually at most of the order of the Young’s modulus, which (for rubber) is less than 10^7 Pa. Thus, in most cases involving elastically soft materials, the viscosity can be considered as independent of the local pressure. In the applications below the nominal pressure is only of the order of ~10^4 Pa and the pressure in the area of real contact of the order of ~10^6 Pa, so that the dependence of the (shear) viscosity on the pressure can be neglected.

It has been observed experimentally [17, 18], and also found in molecular dynamics (MD) simulations [19, 20], that the effective viscosity \( \eta \) (defined by \( \sigma = \eta v/\nu \), where \( \sigma \) is the shear stress, \( \nu \) the separation between the surfaces and \( v \) the relative velocity) of very thin (nanometer thickness) fluid films confined between solid walls at low pressure may be strongly enhanced and to exhibit non-Newtonian properties. In addition, for nanometer wall–wall separations, a finite normal stress is necessary for the squeeze-out, i.e. the ‘fluid’ now behaves as a soft solid and the squeeze-out occurs in a quantized way by removing one monolayer after another with increasing normal stress [21]. In the application below we only study the (average) separation between the walls with micrometer resolution, and in this case the strong increase in the viscosity for very short wall separations becomes irrelevant.

2.3. Roughness on many length scales: effective equations of fluid flow

Equations (2) and (3) describe the fluid flow at the interface between contacting solids with rough surfaces. The surface roughness can be eliminated or integrated out using the renormalization group (RG) procedure. In this procedure one eliminates or integrates out the surface roughness components in steps and obtain a set of RG flow equations describing how the effective fluid equation evolves as more and more of the surface roughness components are eliminated. One can show that, after eliminating all the surface roughness components, the fluid current (given by (2)) takes the form

\[
J = A(\bar{u}) \nabla \bar{p} + B(\bar{u}) v_0
\]

where \( A \) and \( B \) are \( 2 \times 2 \) matrices, and where \( \bar{u}(x, t) \) and \( \bar{p}(x, t) \) are now locally averaged quantities. In general, \( A \) and \( B \) depend also on \( \nabla \bar{p} \) (see [22]), but for the low pressures (and pressure gradients) prevailing in the application presented below, we can neglect this effect.

In the original work of Patir and Cheng it was not proved that the flow current can be written as in (4), but later work has derived (4) rigorously using different procedures, including the RG method [15] and the homogenization procedure described in [23]. The basic assumption necessary is only the separation of length scales: the longest wavelength roughness component should be much smaller than the nominal size of the contact region. The roughness should be random (i.e. the phases of the different surface roughness wavevector component should be uncorrelated), but it does not need to be self-affine fractal, i.e. the surface roughness power spectrum \( C(q) \) can have any form, e.g. it does not need to be a power in the wavevector, as would be the case for a self-affine fractal surface.

If the sliding velocity \( v_0 = 0 \) and if the surface roughness has isotropic statistical properties, then \( A \) is proportional to the unit matrix and is usually written as \( A = -\bar{u}^3 \phi_p(\bar{u})/(12\eta) \). In this case from (3) and (4) we obtain

\[
\frac{\partial \bar{u}}{\partial t} - \nabla \cdot \left( \frac{\bar{u}^3 \phi_p(\bar{u})}{12\eta} \nabla \bar{p} \right) = 0.
\]

In figure 2 we show \( \phi_p(\bar{u}) \) calculated using the Persson contact mechanics and the Bruggeman effective medium theory [15]. The figure shows the dependence of \( \phi_p(\bar{u}) \) on the separation \( \bar{u} \) for the two (copper) surfaces used in the study below. The green curve shows the large \( \bar{u} \) behavior predicted by Tripp [24]:

\[
\phi_p \approx 1 - \frac{3}{2} \frac{\langle h^2 \rangle}{\bar{u}^2}.
\]

where \( \langle h^2 \rangle = h_{rms}^2 \) is the ensemble average of the square of the roughness amplitude. See also [15] for the calculation of higher-order corrections.

2.4. Fluid squeeze-out

Let us squeeze a cylindrical rubber block (height \( d \) and radius \( R \)) against a substrate in a fluid. Assume that we can neglect the macroscopic deformations of the rubber block in response to the (macroscopically) non-uniform fluid pressure. In this case \( \bar{u}(x, t) \) will only depend on time \( t \) and (5) reduces to

\[
\frac{d\bar{u}}{dt} - \bar{u}^3 \frac{\phi_p(\bar{u})}{12\eta} \nabla^2 \bar{p} = 0.
\]

This equation implies that the fluid pressure

\[
\bar{p} = 2\rho_{\text{fluid}} \left( 1 - \frac{\bar{u}^2}{R^2} \right)
\]
where \( r = |x| \) denotes the distance from the cylinder axis, and where we have assumed that the external pressure vanishes. \( p_{\text{fluid}} \) denotes the average fluid pressure in the nominal contact region. Substituting (7) in (6) gives

\[
\frac{d\bar{u}}{dt} \approx \frac{2\bar{u}^2 \phi_p(\bar{u}) p_{\text{fluid}}(t)}{3\eta R^2}.
\]

(8)

If \( p_0 \) is the applied pressure acting on the top surface of the cylinder block, we have

\[
p_{\text{fluid}}(t) = p_0 - p_{\text{cont}}(t),
\]

(9)

where \( p_{\text{cont}} \) is the (locally averaged) asperity contact pressure. We first assume that the pressure \( p_0 \) is so small that for all times \( \bar{u} \gg h_{\text{rms}} \) and in this case \( \phi_p(\bar{u}) \approx 1 \). For \( \bar{u} \gg h_{\text{rms}} \) we also have [9]

\[
p_{\text{cont}} \approx \beta E^* \exp(-\bar{u}/u_0),
\]

(10)

where \( E^* = E/(1-\nu^2) \) (where \( E \) is the Young’s modulus and \( \nu \) the Poisson ratio), and where \( u_0 = h_{\text{rms}}/\alpha \). The parameters \( \alpha \) and \( \beta \) depend on the fractal properties of the rough surface [9].

We note that (10) was derived for the dry condition and is only approximately valid in the present case. Thus, we have assumed that the main effect of the fluid pressure is to reduce the contact pressure \( p_{\text{cont}} = p_0 - p_{\text{fluid}} \). However, the non-uniform (on the asperity length scale) fluid pressure will also deform the surface of the elastic solid non-uniformly, in particular at the length scale determined by the longest wavelength surface roughness components. This topic is discussed in [22, 25] (mainly in the context of mixed lubrication). The elastic deformations induced by the non-uniform (on the asperity length scale) fluid pressure will tend to effectively make \( u(x, t) \) more uniform (as if the surfaces were smoother) which will increase the squeeze-out time and may, in addition to the fluid-induced macroscopic deformations discussed below (resulting from the macroscopic variation of the fluid pressure with the distance \( r \) from the center of the contact region), slow down the squeeze-out. However, because of the low fluid pressure involved in our experiments (see below), this effect is negligible. In particular, if \( Q = (p_0/E)(\lambda/R)/(\lambda/\bar{u}) < 1 \), where \( \lambda \) is the longest wavelength roughness component, then the fluid-induced deformations of the solid walls at the asperity length scale can be neglected [22, 25]. In our case \( p_0/E \approx 0.01, \lambda/R \approx 0.1 \) and \( \lambda/\bar{u} \approx 10 \) so that \( Q \approx 0.01 \).

Using (10) and (9) we get from (8)

\[
\frac{dp_{\text{cont}}}{dt} \approx \frac{2\bar{u}^3 (p_{\text{cont}}(t))}{3\eta R^2 u_0} p_{\text{cont}}(p_0 - p_{\text{cont}}).
\]

(11)

For long times \( p_{\text{cont}} \approx p_0 \) and we can approximate (11) with

\[
\frac{dp_{\text{cont}}}{dt} \approx \frac{2\bar{u}^3 (p_0)}{3\eta R^2 u_0} p_0(p_0 - p_{\text{cont}}).
\]

Integrating this equation gives

\[
p_{\text{cont}}(t) \approx p_0 - \{p_0 - p_{\text{cont}}(0)\} \exp\left(-\frac{(\bar{u}(p_0)/h_{\text{rms}})^3 t}{\tau}\right).\]

(12)
the change in the thickness of the rubber block, the local elastic asperity-induced deformations at the (lower) interface will be determined (to a good approximation) by the Young's modulus $E$.

In section 2.4 we have shown that a flat cylinder surface squeezed against a flat substrate in a (Newtonian) fluid gives rise to a parabolic fluid pressure distribution. This implies that, for a very thin ($d \ll R$) elastic disc, glued to a flat rigid surface and squeezed against another flat surface in a fluid, we expect the bottom surface of the elastic disc to remain nearly flat, and the assumption made in section 2.4 will hold to good accuracy. However, if the rubber block is thick enough ($d > R$) the bottom surface of the block will bend inwards as indicated in figure 4, which will slow down the fluid squeeze-out.

Including the macroscopic deformations of the bottom surface of the cylinder block, in response to the fluid pressure distribution $\hat{p}(x)$, is a very complex (numerical) problem, which we have not been able to solve so far. However, for the case of a rubber sphere (or a half-sphere) the numerical problem of accounting both for the fluid-induced macroscopic deformation, and the asperity interaction, becomes much simpler and we are at present studying this case numerically, and we also plan to perform an experiment for this situation in the near future.

3. Experimental details

We have studied the squeeze-out of a fluid between solids with rough surfaces as shown in figure 5 and in figure 6 schematically. In the experimental set-up a cylindrical silicon rubber block is squeezed against a rough countersurface in the presence of a fluid. The cylindrical body has the height $d = 1, 0.5$ or $0.3$ cm, and the diameter $D = 2R = 3$ cm. The normal load $F_N = 13.8$ N and the fluid viscosity $\eta = 100$ Pa s. The vertical displacement $s(t)$ of the upper surface is registered as a function of time $t$.

![Figure 3](image1.png)

**Figure 3.** A cylindrical rubber block (height $d$ and radius $R$) squeezed against a lubricated substrate (no friction). If $d > R$ the pressure distribution at the interface will be nearly uniform (left) while if $d \ll R$ (right) the pressure distribution will be nearly parabolic. We have assumed that the upper surfaces of the rubber cylinders are glued (no slip) to a flat rigid disc.

![Figure 4](image2.png)

**Figure 4.** The non-uniform hydrodynamic pressure is highest at the center of the contact region and will deform the rubber block as indicated in the figure.

![Figure 5](image3.png)

**Figure 5.** Experimental set-up for studies of fluid squeeze-out between surfaces of elastic solids.

![Figure 6](image4.png)

**Figure 6.** Squeeze-out experimental set-up (schematic). A cylindrical glass or rubber block is squeezed against a substrate with a smooth or rough surface in a fluid. The cylindrical body has the height $d = 1, 0.5$ or $0.3$ cm, and the diameter $D = 2R = 3$ cm. The normal load $F_N = 13.8$ N and the fluid viscosity $\eta = 100$ Pa s. The vertical displacement $s(t)$ of the upper surface is registered as a function of time $t$. 

The normal load is kept constant for all experiments. We have measured the downwards movement of the dead weight as a function of time using a digital gauge with a relative position resolution of $0.5 \, \mu$m. In order to slow down the whole process, we use a very high viscosity silicon oil (Dow Corning 200 Fluid,
we show the power spectrum as a function of the logarithm of the wavevector for the two copper surfaces, 1 and 2, with the root-mean-square roughness 42 µm and 88 µm, respectively.

Figure 7. The logarithm of the surface roughness power spectrum as a function of the logarithm of the wavevector for the two copper surfaces, 1 and 2, with the root-mean-square roughness 42 µm and 88 µm, respectively.

viscosity 100 Pa s) and a relatively low nominal squeezing pressure (about 10^4 Pa). In the different configurations we either squeeze an elastic silicon rubber block, or a rigid glass block, against smooth (glass) or rough (copper) surfaces in order to test different aspects of the squeeze-out. The rubber blocks have the radius \( R = 1.5 \) cm and height \( d = 1 \), 0.5 and 0.3 cm. We use a silicone elastomer (PDMS) prepared using a two-component kit (Sylgard 184) purchased from Dow Corning (Midland, MI). This kit consists of a base (vinyl-terminated polydimethylsiloxane) and a curing agent (methylhydroxiloxane–dimethylsiloxane copolymer) with a suitable catalyst. From these two components we prepared a mixture 10:1 (base/cross linker) in weight. The mixture was degassed to remove the trapped air induced by stirring from the mixing process and then poured into cylindrical casts. The bottom of these casts was made from glass to obtain smooth surfaces (negligible roughness). The samples were cured in an oven at 80 °C for over 12 h. The rough copper surfaces were prepared by pressing sandpaper surfaces against flat and plastically soft copper surfaces using a hydraulic press. Using sandpaper with different grit sizes, and repeating the procedure many times, resulted in (nearly) randomly rough surfaces suitable for our experiment.

The silicon block was placed in the high viscosity fluid with some distance to the rough surface. In order to avoid kinetic (inertia) effects, the initial separation was selected to be very small. The nominal force was applied by dropping the dead weight with the rubber block attached to it. The displacement of the dead weight from its starting position was measured as a function of time.

The surface topography \( z = h(x) \) of the two copper surfaces 1 and 2 used in our study has been measured using white-light interferometry. White light is focused on the substrate and the topography is determined by analyzing the scattered light and its wavelength. Using this optical method, the roughness of the soft copper samples can be studied without destroying the surface. The vertical resolution of the sensor is 20 nm while the resolution in the xy plane is 2 µm. In figure 7 we show the power spectrum \( C(q) \) of the two copper surfaces 1 and 2 defined by [7]

\[
C(q) = \frac{1}{(2\pi)^2} \int d^2 x \langle h(x)h(0) \rangle e^{-iq \cdot x}
\]

where \( \langle \cdots \rangle \) stands for the ensemble average.

The area of real contact (at the nominal squeezing pressure \( \approx 2 \times 10^4 \) Pa) as a function of the magnification \( \zeta \) is shown in figure 8. Note that the area of real contact (i.e. the contact area (projected on the xy plane) at the highest magnification \( \zeta_1 \) or wavevector \( q_1 = q_0 \zeta_1 \)) is rather similar in both cases (equal to \( A = 0.016A_0 \) and 0.013\( A_0 \) for surfaces 1 and 2, respectively) in spite of the rather large difference in the rms roughness values (\( h_{rms} = 42 \) and 88 µm, respectively). This is due to the fact that the rms roughness is dominated by the longest wavelength roughness components, while the area of real contact is strongly influenced by the short wavelength roughness components, which are very similar on both surfaces (see figure 7 for large wavevector). The small contact pressure results in a relative large (average) separation between the surfaces, \( \bar{u} \approx 1.4h_{rms} \), for both surfaces (as calculated using the theory developed in [9, 10]).

4. Comparison of theory with experiment

In figure 9 we show the surface separation as a function of the logarithm of time when a glass and a PDMS cylindrical block are squeezed against a flat glass substrate in a silicon oil. Also shown is the theoretical prediction (lower curve). The cylinder is \( d = 0.5 \) cm thick and has the diameter \( D = 2R = 3 \) cm. As expected, the theoretical result agrees almost perfectly with the experimental results for the glass cylinder (no fitting parameters), but for the rubber block the (average) separation is larger and the squeeze-out slower. We attribute this to temporarily trapped fluid resulting from the upward bending (before contact with the substrate) of the bottom surface of the rubber block, as indicated in figure 4. We define the ‘trapped’ fluid volume \( \Delta V \) as the fluid volume between the
rubber is smaller for the thin rubber disc (see section fluid-pressure-induced curvature of the bottom surface of the thinner rubber disc. This is indeed expected since the agreement between the theory and experiment is much better. 

The surface separation as a function of the logarithm of time when a glass and a PDMS cylindrical block is squeezed against a flat glass substrate in a silicon oil. Also shown is the theoretical prediction (lower curve). The cylindrical body has the height \( d = 0.5 \text{ cm} \) and diameter \( D = 2R = 3 \text{ cm} \). The normal load \( F_N = 13.8 \text{ N} \) and the fluid viscosity \( \eta = 100 \text{ Pa s} \).

Figure 9. The surface separation as a function of the logarithm of time when a glass and a PDMS cylindrical block is squeezed against a flat glass substrate in a silicon oil. Also shown is the theoretical prediction (lower curve). The cylindrical body has the height \( d = 0.5 \text{ cm} \) and diameter \( D = 2R = 3 \text{ cm} \). The normal load \( F_N = 13.8 \text{ N} \) and the fluid viscosity \( \eta = 100 \text{ Pa s} \).

The surface separation as a function of the logarithm of time when PDMS cylindrical blocks with thicknesses 1 cm and 0.3 cm are squeezed against a rough copper surface 1 (root-mean-square roughness \( h_{\text{rms}} = 42 \text{ \mu m} \)) in a silicon oil. Also shown is the theoretical prediction for a flat substrate (dashed curve) and for copper surface 1 (lower solid curve). The cylindrical body has the height \( D = 10 \text{ mm} \), the radius \( R = 5 \text{ cm} \), the thickness \( d = 2.5 \text{ mm} \), the normal load \( F_N = 100 \text{ N} \) and the fluid viscosity \( \eta = 100 \text{ Pa s} \).

Figure 10. (a) The surface separation as a function of the logarithm of time when PDMS cylindrical blocks with thickness 1 cm and 0.33 cm are squeezed against a rough copper surface 1 (root-mean-square roughness \( h_{\text{rms}} = 42 \text{ \mu m} \)) in a silicon oil. Also shown is the theoretical prediction for a flat substrate (dashed curve) and for copper surface 1 (lower solid curve). (b) The same as in (a) but for a more narrow time interval. The thin solid line is the calculated squeeze-out when the pressure flow factor \( \phi_p(\bar{u}) = 1 \).

contact with the substrate, the fluid-pressure-induced bending of the bottom surface of the block is the same in both cases (giving overlapping curves for \( t < 300 \text{ s} \)).

In figure 12 we show the surface separation as a function of the logarithm of time when the \( d = 0.5 \text{ cm} \) thick PDMS cylindrical block is squeezed against the (rough) copper surface 2 (root-mean-square roughness \( h_{\text{rms}} = 88 \text{ \mu m} \)) in a silicon oil. Also shown is the theoretical prediction (lower solid curve). Again, the bending of the bottom surface of the rubber block results in a slower squeeze-out than predicted theoretically assuming a (macroscopically) flat bottom surface of the rubber block.

We are convinced that the discrepancy between our calculations and the experimental data is due to the fluid-induced bending of the bottom surface of the block. As far as we know, this effect has not been studied experimentally before for the cylinder-block geometry. However, it is very well known (both from experiments and from elastohydrodynamic calculations) that, for an elastic sphere squeezed against a flat in a fluid, the fluid pressure is highest at the center of the ‘contact area’ and bends the
Figure 11. The surface separation as a function of the logarithm of time when a 0.5 cm thick PDMS cylindrical block is squeezed against a flat glass substrate (lower curve) and against the (rough) copper surface 1 (root-mean-square roughness $h_{\text{rms}} = 42 \, \mu\text{m}$) in a silicon oil.

Figure 12. The surface separation as a function of the logarithm of time when a 0.5 cm thick PDMS cylindrical block is squeezed against the rough copper surface 2 (root-mean-square roughness $h_{\text{rms}} = 88 \, \mu\text{m}$) in a silicon oil. Also shown is the theoretical prediction (lower solid curve). The dashed curve is the theoretical prediction for a flat (no surface roughness) substrate.

bottom surface of the sphere inwards so that the solid–solid separation is largest at the center (where the separation would be smallest if no deformation of the ball would occur) [5]. For the cylinder-block case (if the thickness is of the order of or larger than the diameter of the cylinder) this bending will be even larger.

5. Summary and conclusion

In this paper we have studied the fluid squeeze-out from the interface between an elastic block with a flat surface and a randomly rough surface of a rigid solid. We have calculated the (average) interfacial separation as a function of time by considering the fluid flow using a contact mechanics theory in combination with thin-film hydrodynamics with flow factors (which are functions of the (local) interfacial separation) obtained using a recently developed theory. We have shown the importance of the large length-scale elastic deformations on the squeeze-out.

The theoretical results have been compared to experimental results. The experiment was performed by squeezing cylindrical rubber blocks with different heights $d$ against rough cooper surfaces in the presence of a high viscosity fluid (silicone oil). Changing the height $d$ of the rubber block, and also performing additional experiments with flat against flat surfaces, with combinations of rigid–rigid and elastic–rigid, we could show the importance of both the large length scale and asperity-induced elastic deformation on the squeeze-out. In particular, large length-scale deformations of the bottom surface of the rubber block resulted in (temporary) trapped fluid between the elastic solid and the rigid countersurface, which drastically slowed down the squeeze-out. This effect is smallest for the thinnest rubber block, in which case we find good agreement between the theory (where we have neglected the large length-scale deformations of the rubber block) and the experiments. Another mechanism which drastically slows down the squeeze-out occurs at much higher nominal pressure (or load) than used in the present experiment. This is due to sealed-off fluid in the nominal contact region during contact formation. This effect occurs when the area of real contact approaches $\approx 0.5 A_0$, where the area of real contact percolates, resulting in sealed-off regions of fluid, which may disappear only extremely slowly, e.g. by diffusion of the fluid into the rubber. This effect was discussed in [12] and seems to be of importance in many applications involving high contact pressures, e.g. it may result in a static (or start-up) friction force which slowly increases with time even after very long times (say, one year).

The squeeze-out of a fluid from the interfacial region between elastic solids with rough surfaces is very important in many technical applications (e.g. a tire rolling on a wet road, wipers and dynamic seals), and the results presented in this paper contribute to this important subject.

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