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Rolling Friction: Comparison of Analytical Theory with Exact Numerical Results

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Abstract We study the contact mechanics of a smooth hard cylinder rolling on a flat surface of a linear viscoelastic solid. Using the measured viscoelastic modulus of unfilled and filled (with carbon black) nitrile rubber, we compare numerically exact results for the rolling friction with the prediction of a simple analytical theory. For the unfilled rubber, the two theories agree perfectly while some small difference exists for the filled rubber. The rolling friction coefficient depends nonlinearly on the normal load and the rolling velocity.

Keywords Rubber friction · Rolling friction · Polymer · Viscoelasticity

1 Introduction

Rubber friction is a topic of huge practical importance, e.g., for tires, rubber seals, wiper blades, conveyor belts, and syringes [1–16]. Many experiments have been performed with a hard sphere or cylinder rolling on a flat rubber substrate [17–20]. Nearly the same friction force may be observed during sliding as during rolling, assuming that the interface is lubricated and that the sliding velocity and fluid viscosity are such that a thin lubrication film is formed with a thickness much smaller than the indentation depth of the ball,

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but larger than the amplitude of the roughness on the surfaces [17]. The results of rolling friction experiments have often been analyzed using a very simple model of Greenwood and Tabor [17], which, however, contains a (unknown) factor α , representing the fraction of the input elastic energy lost as a result of the internal friction of the rubber.

Several (analytical) theoretical studies have been presented for the rolling and sliding friction involving viscoelastic solids, which, however, are based on a simplified description of the rubber rheology, i.e., on a single term Prony series. The case of cylinder contact has been analytically treated by Hunter [21] and Goriacheva [22], and the approach lately approximately extended to the case of three-dimensional contact by enforcing a two-dimensional reduction in the problem [23]. However, the relaxation spectrum of real polymers is usually very wide and the rolling friction of real rubber compounds cannot be predicted using those existing exact theories. This limitation also applies to recent numerical models based on the efficient boundary element formulation of the contact problem [24], where the single relaxation time rheology is adopted.

Recently, one of us (BNJP) has developed a Fourier series-based approach for the prediction of rolling friction for real rubber materials [25]. This approach, accurate to linear order in the loss tangent, is able to predict quantitatively reliable results, as shown for both line [25] (see also in the following) and point contacts [26], with negligible computational effort. The theory can be applied to solids with arbitrary viscoelastic modulus. For a cylinder rolling on a viscoelastic solid characterized by a single relaxation time, Hunter [21] has obtained an exact result for the rolling friction, and the analytical theory [25] agrees very well with the result of Hunter. Unfortunately, the approach in Ref. [25] does not allow us to determine the traction and the displacement fields occurring at the contact.



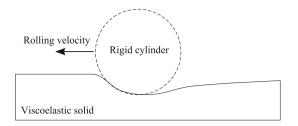


Fig. 1 Smooth rigid cylinder in steady rolling contact with a viscoelastic solid with nominally flat surface. Schematic

Full numerical solutions of the viscoelastic contact problem, of interest for many tribological applications, have only very recently been presented for smooth and rough contacts, see e.g., Ref. [26]. This should not surprise at all, due to the numerical complexity of the underlying problem, which is even more pronounced in case of real (multiscale) rough contacting surfaces. As a result, the hysteretic rubber friction coming from local sliding contacts can be, nowadays, only effectively calculated based on homogenized contact mechanics theories [3].

In this study, we compare the analytical theory of Ref. [25] with exact numerical results. In particular, in Sect. 2, we describe our numerical approach; in Sect. 3, we briefly describe the analytical approach of Ref. [25] as applied to cylinder on flat. Section 4 presents the numerical comparison of the two models. Section 5 contains the summary and conclusion.

2 Rolling Friction: Numerical Theory

We consider the line contact case with a rigid cylinder, with perfectly smooth surface, rolling on an isotropic viscoelastic half-space, see Fig. 1. In order to determine the surface deformation of the viscoelastic half-space, we assume that quasi-static bulk dynamics holds, i.e., inertia effects are neglected. If a stress $\sigma_z(\mathbf{x},t)$ acts on the surface of a semi-infinite viscoelastic solid, it will result in a normal surface displacement $u_z(\mathbf{x},t)$. The latter is linearly related to $\sigma_z(\mathbf{x},t)$ via an equation which is particular simple when the spatial and time coordinates are Fourier transformed [3]:

$$u_{\mathbf{z}}(\mathbf{q},\omega) = M_{\mathbf{z}\mathbf{z}}(\mathbf{q},\omega)\sigma_{\mathbf{z}}(\mathbf{q},\omega),$$
 (1)

where

$$M_{zz}(\mathbf{q},\omega) = -\frac{2(1-v^2)}{qE(\omega)},\tag{2}$$

and where $E(\omega)$ is the frequency-dependent (complex) Young's modulus. In what follows, we will assume that the Poisson's ratio $v(\omega)$ is independent of frequency. This

assumption is a good approximation in most cases (see e.g., Ref. [25]), and it means that the dynamics of relaxation in orthogonal directions proceeds equally in time. If a rigid cylinder is sliding in steady state on a semi-infinite rubber solid, we have

$$\sigma_{\mathbf{z}}(\mathbf{x},t) = \sigma_{\mathbf{z}}(\mathbf{x} - \mathbf{v}t)$$

which gives

$$\sigma_{z}(\mathbf{q},\omega) = \delta(\mathbf{q} \cdot \mathbf{v} - \omega)\sigma_{z}(\mathbf{q}). \tag{3}$$

For the cylinder geometry with the cylinder axis along the y-direction, $\sigma_z(\mathbf{x})$ is independent of y, and we write $\sigma_z(\mathbf{x}) = -p(x)$ so that

$$\sigma_{\mathbf{z}}(\mathbf{q}) = -p(q_{\mathbf{x}})\delta(q_{\mathbf{y}}). \tag{4}$$

Using (1)-(4) gives

$$u_{\rm z}(\mathbf{q},\omega) = \frac{2(1-v^2)}{|q_{\rm x}|E(\omega)}p(q_{\rm x})\delta(q_{\rm y})\delta(q_{\rm x}v_{\rm x}-\omega).$$

Thus, we get:

$$u_{z}(x,t) = \int d^{2}q d\omega \ u_{z}(\mathbf{q},\omega) e^{i(\mathbf{q}\cdot\mathbf{x}-\omega t)}$$

$$= \int d\omega \frac{2(1-v^{2})}{|\omega|E(\omega)} p(\omega/v_{x}) e^{i\omega(x/v_{x}-t)}. \tag{5}$$

Now because of causality, one can write

$$\frac{1}{E(\omega)} = \frac{1}{E(\infty)} + \int_{0}^{\infty} d\tau \frac{H(\tau)}{1 - i\omega\tau},$$
 (6)

where the spectral density $H(\tau)$ is a positive real function of the rubber relaxation time τ . Note that we can also write

$$\frac{1}{E(\omega)} = \frac{1}{E(\infty)} \left[1 + \int_{0}^{\infty} dt \ \Gamma(t) e^{i\omega t} \right], \tag{7}$$

where

$$\Gamma(t) = E(\infty) \int_{0}^{\infty} d\tau \, \frac{H(\tau)}{\tau} e^{-t/\tau}.$$

We write

$$p(q_{x}) = \frac{1}{2\pi} \int dx' \ p(x') e^{-iq_{x}x'}.$$
 (8)

Substituting (7) and (8) in (5) gives

$$u_{z}(x,t) = \frac{(1-v^{2})}{\pi E(\infty)} \int dx' \int d\omega \frac{1}{|\omega|} p(x')$$

$$\times \left[e^{i\alpha\omega/v_{x}} + \int_{0}^{t} dt' \ \Gamma(t') e^{i\beta(t')\omega/v_{x}} \right], \tag{9}$$



where $\alpha = x - x' - v_x t$ and $\beta(t') = x - x' - v_x (t - t')$. Now consider the integral

$$J=\int\limits_{-\infty}^{\infty}\,\mathrm{d}\omega\,rac{1}{|\omega|}e^{ilpha\omega/v_x}.$$

We write $\omega/v_x = q_x$ and get

$$J = 2 \int_{0}^{\infty} \mathrm{d}q_{\mathrm{x}} \, \frac{\cos(\alpha q_{\mathrm{x}})}{q_{\mathrm{x}}}.$$

This integral diverges. The problem arises from the small wavevector region $q_x \to 0$ and reflects the infinite size of the system. For a finite (but very large) solid with linear dimension L, we must replace the lower integration limit with π/L , so that

$$J = 2 \int_{\pi/L}^{\infty} \mathrm{d}q_{\mathrm{x}} \frac{\cos(\alpha q_{\mathrm{x}})}{q_{\mathrm{x}}}.$$

Let us denote $|\alpha|q_{\rm x}=\xi$ so that

$$J = 2 \int_{|\alpha|\pi/L}^{\infty} d\xi \, \frac{\cos \xi}{\xi}.$$

For large (but fixed) L, this will give

 $J \approx \text{const.} - 2\log|\alpha|$.

Using this result in (9) gives

$$u_{z}(x,t) = u_{0} - \frac{2(1-v^{2})}{\pi E(\infty)} \int dx' \ p(x') \times \left[\log|x - x' - v_{x}t| + \int_{0}^{t} dt' \ \Gamma(t') \log|x - x' - v_{x}(t - t')|, \right]$$

where u_0 is a constant. In what follows, we study the system in a coordinate system fixed to the cylinder which correspond to replacing $x \to x + v_x t$. We also denote the velocity of the cylinder with $v_x = v_0$ and the high-frequency modulus $E(\infty)$ with E_∞ . We consider long times $t \to \infty$ where the sliding is steady state, where u(x,t) = u(x) (in the moving reference frame) is time independent. With these assumptions, we get

$$u_{z}(x) = u_{0} - \frac{2(1 - v^{2})}{\pi E_{\infty}} \int dx' \ p(x')$$

$$\times \left[\log|x - x'| + \int_{0}^{\infty} dt' \ \Gamma(t') \log|(x - x' - v_{0}t')| \right].$$
(10)

Eq. (10) is based on a Green's function formulation of the bulk viscoelastic deformation problem, similarly to Ref. [26]. The derivation above is for a homogeneous viscoelastic half-space, but can be easily generalized to a layered material, which introduce in (2) an additional factor $S(q,\omega)$ which is known analytically (see e.g., Ref. [27]). In this work, differently from what reported in Ref. [26], the line contact condition allows us to derive an exact discrete formulation of Eq. (10) [28], hence avoiding numerical issues coming from its integration over a wide range of relaxation times describing real rubbers.

The rolling friction coefficient:

$$\mu_{\rm R} = -\frac{1}{f_{\rm N}} \int dx \ p(x) h'(x),$$

where $h(x) = \text{const.} - (R^2 - x^2)^{1/2} \approx \text{const.} + x^2/2R$ is the cylinder surface height.

The constant u_0 in (10) can be implicitly calculated with the load balance equation

$$\int_{-\infty}^{\infty} dx' \ p(x') = f_{\rm N},\tag{11}$$

where $f_N = F_N/L_y$ is the cylinder load per unit length. Equations (10) and (11) together with the relation between the interfacial separation and the nominal pressure given by the Persson's contact mechanics theory [29] have to be solved to determine the unknown variables p(x), u(x) and u_0 . Here, we use such small surface roughness that it has a negligible influence on the rolling friction result, which is therefore essentially exact.

3 Rolling Friction: Analytical model

If a solid is exposed to a perpendicular surface stress or pressure $p(\mathbf{x},t) = p(\mathbf{x} - \mathbf{v}t)$, a tangential (friction) force will act on the solid given by [25]

$$F_{\rm f} = \frac{2(2\pi)^2}{v} \int d^2q \, \frac{\omega}{q} \text{Im} \frac{1}{E_{\rm eff}(\omega)} |p(\mathbf{q})|^2, \tag{12}$$

where $\omega = \mathbf{q} \cdot \mathbf{v}$ and $E_{\rm eff} = E/(1-v^2)$. In principle, v depends on frequency but the factor $1/(1-v^2)$ varies from $4/3 \approx 1.33$ for v = 0.5 (rubbery region) to ≈ 1.19 for v = 0.4 (glassy region), and we can neglect the weak dependence on frequency. Within the assumptions given above, Eq. (12) is exact. Note that even if we use $p(\mathbf{q})$ calculated to zero order in $\tan \delta$, the friction force (12) will be correct to linear order in $\tan \delta$. We also note that the present approach is very general and flexible. For example, instead of a semi-infinite solid, the theory is easily generalized to layered materials, e.g., a thin viscoelastic film on a hard flat substrate.



$$p(x) = \frac{2f_{\rm N}}{\pi a} \left[1 - \left(\frac{x}{a}\right)^2 \right]^{1/2},\tag{13}$$

where $f_{\rm N}=F_{\rm N}/L_{\rm y}$ the load per unit length of the cylinder. The half-width of the contact area in the Hertz contact theory is

$$a = \left(\frac{4f_{\rm N}R}{\pi E_{\rm eff}}\right)^{1/2},\tag{14}$$

where we take $E_{\rm eff}$ to be $|E_{\rm eff}(\omega)|$ with $\omega = \mathbf{q} \cdot \mathbf{v}$. Substituting the Fourier transform of (13) in (12) and dividing by $F_{\rm N}$ gives the friction coefficient [25]:

$$\mu_{\rm R} = \frac{8f_{\rm N}}{\pi} \int_0^\infty dq_{\rm x} \, \, {\rm Im} \frac{1}{E_{\rm eff}(q_{\rm x}v)(aq_{\rm x})^2} J_1^2(q_{\rm x}a), \tag{15}$$

where $J_1(x)$ is the Bessel function of the first kind.

At this point, we empathize that (12) is basically exact, and if one could calculate the contact pressure p(x) [or rather the Fourier transform $p(q_x)$] exactly, then (12) should give the exact result, e.g., the result of Hunter for the cylinder case when assuming the simple viscoelastic modulus consisting of a single term Prony series. Note that the contact pressure will not be symmetric around the midpoint when the dissipation in the rubber is included in the analysis. Still, to linear order in $\tan \delta$, one can neglect this effect [since (12) is explicitly already linear in $\tan \delta$], and the analysis below and in Ref. [25] shows that this is a remarkable accurate approximation, which is an interesting result in its own right, and makes it possible to apply the theory to a wide set of problems.

4 Numerical Results

Here, we compare the results of the numerical procedure developed in Sect. 2 with the predictions of the rolling friction model [25] presented in Sect. 3. Due to the extremely reduced computational requirements of the rolling friction model presented in Sect. 3, with respect to any classical numerical formulation of the viscoelastic contact

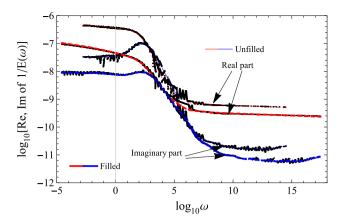


Fig. 2 Real and imaginary part of the complex function $1/E(\omega)$ as measured (*dots*) and fitted by Prony series (*dashed blue and red curves*). For filled (*thick line*) and unfilled rubber. In $\log_{10} - \log_{10}$, with $1/E(\omega)$ in Pa^{-1} and ω in s^{-1} . For the reference temperature, T = 20 °C (Color figure online)

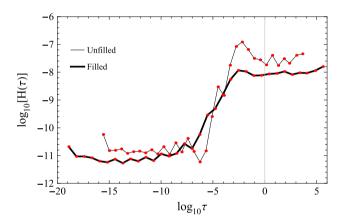


Fig. 3 Spectral density $H(\tau)$ as a function of the relaxation time τ on a $\log_{10}-\log_{10}$ scale. For both filled and unfilled rubber, $H(\tau)$ is Pa^{-1} with τ in s. For the reference temperature, $T=20~\mathrm{^{\circ}C}$

problem¹, it is of particular interest to quantify the accuracy of the approach presented in Sect. 3 in the case of real rubber rheologies. Indeed, while in the case of (unrealistic) simple rheologies, i.e., characterized by a single relaxation time, the theory is very accurate (see Refs. [25, 26]), there is no proof in the literature of its accuracy for a rheology described by several relaxation times.

We have measured the viscoelastic modulus of an unfilled and filled (with carbon black) nitrile rubber. In Fig. 2, we show for both materials the real and imaginary part of the complex function $1/E(\omega)$ as measured (dots) and



¹ It can be easily shown that the discrete formulation of the viscoelastic line contact kernel presents an analytical formulation as it occurs for the elastic case. This largely simplifies the viscoelastic integration.

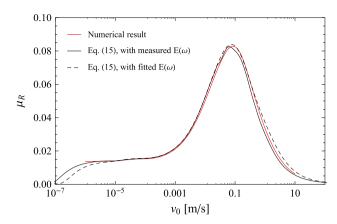


Fig. 4 Rolling friction as a function of the rolling velocity for an unfilled rubber compound, similar to the one described in Fig. 2. For $T=20~^{\circ}C$, R=1~mm and $f_{\rm N}=117~\text{N/m}$

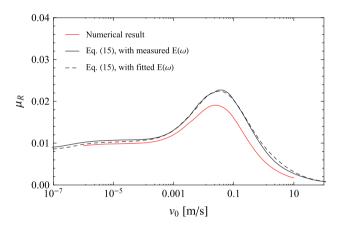


Fig. 5 Rolling friction as a function of the rolling velocity for an filled rubber compound, similar to the one described in Fig. 2. For T = 20 °C, R = 1 mm and $f_N = 117$ N/m

fitted by Prony series (dashed blue and red curves) obtained by approximation (6) with:

$$\frac{1}{E(\omega)} \approx \frac{1}{E(\infty)} + \sum_{n=1}^{N} \frac{H(\tau_n)}{1 - i\omega\tau_n},$$

where N is the number of relaxation times. In Fig. 3, we show the spectral density $H(\tau)$ for both the filled and unfilled rubber. Note that the distribution of relaxation times is very wide and cannot be characterized by a simple single relaxation time as often assumed.

In Figs. 4 and 5, we report, respectively, the rolling friction as a function of the sliding velocity for the unfilled and filled rubber compound. In particular, the red curves are calculated from the numerical theory developed in Sect. 2, whereas the black solid and dashed curves represent, respectively, the rolling friction as calculated from the

analytical theory (Sect. 3) with the measured complex elastic modulus $E(\omega)$ data and from $E(\omega)$ as calculated from the Prony series fitting of the $1/E(\omega)$ measurements. The latter set of data is also used in the numerical exact theory (red curve).

For the unfilled rubber compound, there is a remarkable agreement between the analytical theory of Sect. 3 and the numerical theory of Sect. 2. For the filled rubber, the match is less accurate, and we suspect that the fitting algorithm has been not perfectly successful to accurately determine the rubber relaxed elastic modulus, which is used in our model in the filled case. Due to the sensibility to such a parameter, this could generate a deviation from the results calculated with the Persson's approach [25]. Indeed, it is well know that rubber with filler particles (e.g., carbon black or silica particles) exhibits a strongly nonlinear viscoelastic response, undergoing strain softening already at a strain of order ~ 1 %. A causal linear response function can be written as a Prony series, but this is not the case when the response is nonlinear (see e.g., [30]). Hence, the analytical model, which makes use of the measured frequency space characterization of the rubber, is expected to provide an accurate rolling friction prediction, almost independently the filling state of the rubber. Differently, we suspect that the accuracy of the numerical model, as well as of all existing models based on the space of time formulation of viscoelasticity, can suffer to some extent of the nonlinearity of the rubber rheology when using Prony series fitting.

For the unfilled sample, in Fig. 6 we show the contact pressure field for some representative values of sliding velocity. The different stages of viscoelastic response are highlighted. The strong asymmetry of the contact pressure fields with respect to the Hertzian shape, prevailing in the rubber transition regime, may suggest a certain restriction of accuracy of Eq. (15) in predicting rolling friction. It would be therefore very interesting to compare $|p(q)|^2$ $|f_N J_1(qa)/(\pi qa)|^2$ with respect to the numerically exact results. In particular, in Fig. 7, we show the square of the Fourier transform of the contact pressure (unfilled sample) for representative values of sliding velocities belonging to the rubber transition regime, and corresponding to the pressure curves shown in Fig. 6. Note that all $|p(q)|^2$ curves converge to the same value for $q \approx 0$, as expected since $p(q=0) = f_N/(2\pi)$. Most interestingly, despite the strong asymmetrical shape of the pressure curves in the transition regime (Fig. 6), in the Fourier space, the inaccuracy resulting from such asymmetry is notably reduced, as expected by previous arguments.

Note that the rolling friction exhibits a very strong temperature dependence (not shown), given by the WLF shift factor a_T : reducing the temperature by 10 °C typically



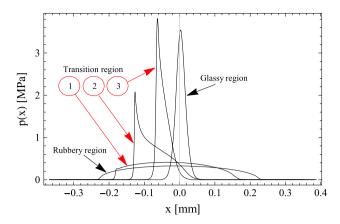


Fig. 6 Contact pressure field for (1) $v_0 = 3.0 \text{ mm/s}$, (2) $v_0 = 77 \text{ mm/s}$, (3) $v_0 = 0.39 \text{ m/s}$. For the unfilled rubber, and for the same parameters of Fig. 4

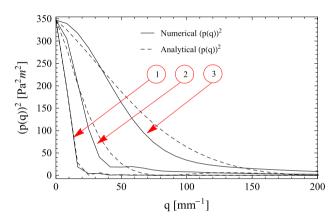
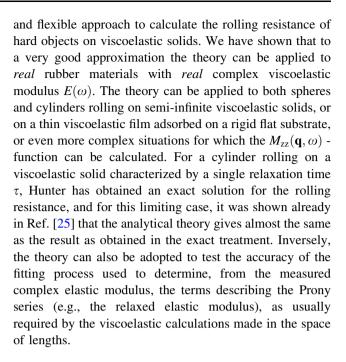


Fig. 7 Square of the Fourier transform of the contact pressure for the unfilled sample. For representative values of sliding velocities belonging to the rubber transition regime and corresponding to the pressure curves shown in Fig. 6

results in a shift of the rolling friction curve toward lower velocities by approximately one decade. At low frequencies (or low rolling velocities), the unfilled rubber is elastically much softer than the filled rubber. Thus, the half-width a_0 of the static Hertz contact region at low rolling velocities is much larger for the unfilled rubber than for the filled rubber. At high rolling velocities, these values becomes much smaller owing to the stiffening of the effective elastic modulus at high frequencies.

5 Summary and Conclusion

We have compared the prediction of an exact numerical study for the rolling friction of a hard cylinder on a smooth flat surface of a viscoelastic solid, with the prediction of an analytical theory. The analytical theory is a very general



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