

Data-driven prediction of the equivalent sand-grain height in rough-wall turbulent flows

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This paper investigates a long-standing question about the effect of surface roughness on turbulent flow: What is the equivalent roughness sand-grain height for a given roughness topography? Deep neural network (DNN) and Gaussian process regression (GPR) machine learning approaches are used to develop a high-fidelity prediction approach of the Nikuradse equivalent sand-grain height k_s for turbulent flows over a wide variety of different rough surfaces. To this end, 45 surface geometries were generated and the flow over them simulated at $Re_\tau = 1000$ using direct numerical simulations. These surface geometries differed significantly in moments of surface height fluctuations, effective slope, average inclination, porosity and degree of randomness. Thirty of these surfaces were considered fully rough, and they were supplemented with experimental data for fully rough flows over 15 more surfaces available from previous studies. The DNN and GPR methods predicted k_s with an average error of less than 10 % and a maximum error of less than 30 %, which appears to be significantly more accurate than existing prediction formulae. They also identified the surface porosity and the effective slope of roughness in the spanwise direction as important factors in drag prediction.

Key words: turbulence modelling

1. Introduction

At sufficiently high Reynolds numbers all surfaces are hydrodynamically rough, as is almost always the case for flows past the surfaces of naval vessels. Reviews of roughness effects on wall-bounded turbulent flows are provided by Raupach, Antonia & Rajagopalan (1991) and Jiménez (2004). The most important effect of surface roughness

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in engineering applications is an increase in the hydrodynamic drag (Flack 2018), which is due predominantly to the pressure drag generated by the small-scale recirculation regions associated with individual roughness protuberances.

For the foreseeable future, the most practical approach to making predictive flow calculations for many realistic applications is to use engineering one-point closures of turbulence, such as two-equation turbulent eddy-viscosity models to the Reynolds-averaged Navier–Stokes equations. Existing rough-wall corrections to this type of closure typically model the increase in hydrodynamic drag on a single length scale – the equivalent sand-grain height (Nikuradse 1933) k_s – without physically resolving the surface or changing the governing equations. In the fully rough flow regime, where the wall friction depends on the roughness alone and is independent of the Reynolds number, k_s was observed to quantify the increase in hydrodynamic drag through the empirical relation with the roughness function (defined as the offset of the log–linear velocity profile of a rough-wall flow relative to that of a smooth-wall one),

$$\Delta U^+ = \frac{1}{\kappa} \ln k_s^+ - 3.5, \quad (1.1)$$

where $\kappa = 0.41$ is the von Kármán constant and + represents normalization in wall units.

A universal length scale (e.g. k_s in Nikuradse’s relation, or ϵ in the Moody diagram Moody 1944) that can predict accurately the surface drag coefficient is not known *a priori* and does not appear to be equivalent to any single geometrical length scale, such as an average or a root mean square (r.m.s.) of roughness height (Flack 2018). It is also well-established that k_s can depend on many geometrical parameters such as the effective slope (Napoli, Armenio & De Marchis 2008; Yuan & Piomelli 2014a) and the skewness of the roughness height distribution (Flack & Schultz 2010). Readers are referred to Flack & Schultz (2010) and Bons (2002) for extensive reviews on this topic. Empirical expressions for k_s based on a small number of geometrical roughness parameters include, among others,

$$k_s = c_1 k_{avg} (\alpha_{rms}^2 + c_2 \alpha_{rms}), \quad k_s = c_1 k_{avg} A_s^{c_2} \quad \text{and} \quad k_s = c_1 k_{rms} (1 + S_k)^{c_2}, \quad (1.2a-c)$$

proposed by Bons *et al.* (2001), van Rij, Belnap & Ligrani (2002) and Flack & Schultz (2010), respectively. Here k_{avg} is the average height, α is the local streamwise slope angle and $A_s = (S/S_f)(S_f/S_s)^{-1.6}$ (where S , S_f , S_s are, respectively, the platform area, the total frontal area and the total windward wetted area of the roughness) while k_{rms} and S_k are the r.m.s. and skewness of the roughness height fluctuations and c_1 and c_2 are constants.

The hydrodynamic length scale k_s appears to be correlated with different sets of geometrical parameters for each type of rough surface and no universal correlation currently exists for flow over surfaces of arbitrary roughness. For example, for synthetic roughness comprising closely packed pyramids (Schultz & Flack 2009) and random sinusoidal waves (Napoli *et al.* 2008), it has been shown that k_s scales on the effective slope when the surface slope is gentle (i.e. within the ‘waviness’ regime), whereas the skewness and r.m.s. height, but not slope magnitude, become important when the slope is steeper (i.e. within the ‘roughness’ regime). The boundary between these two regimes has been shown to be surface dependent (Yuan & Piomelli 2014a).

Some more recent studies of k_s correlations are summarized below. Thakkar, Busse & Sandham (2017) carried out direct numerical simulation (DNS) of transitionally rough turbulent flows for different irregular roughness topographies. They found that the roughness function is influenced by solidity, skewness, the streamwise correlation

length scale and the r.m.s. of roughness height. Flack, Schultz & Barros (2020) performed several experiments to systematically investigate the effects of the skewness and amplitude of roughness height on the skin friction. They found that the r.m.s. and skewness of roughness height fluctuations are important scaling parameters for the prediction of roughness function; however, the surfaces with positive, negative and zero skewness values needed different correlations. Also, Chan *et al.* (2015) simulated turbulent pipe flows over sinusoidal roughness geometries and confirmed strong dependence of roughness function on the average height and streamwise effective slope.

In previous studies, the small number of roughness parameters used to devise k_s correlations tended to limit their application to a narrow range of surface roughness. Since it appears that many geometrical parameters, such as porosity, moments of roughness height (e.g. r.m.s., skewness and kurtosis), effective slope and surface inclination angle might affect k_s , it is useful to employ a data science approach suited to modelling large multivariate/multioutput systems.

Specifically, we use machine learning (ML) to explore k_s -prediction approaches that depend on a large set of surface-topographical parameters, with the expectation that the resulting models may be applied accurately to a wider range of surfaces. Since the prediction of k_s from surface topography is essentially a labelled regression problem, supervised ML operations were performed using deep neural networks (DNN) and Gaussian process regressions (GPR). Both methods are explained thoroughly in §3. Readers are referred to the monograph by Rasmussen & Williams (2006) and the review provided by LeCun, Bengio & Hinton (2015) for detailed descriptions of these methods.

An initial ensemble of 60 sets of data on k_s as a function of topographical parameters – 45 DNS results and 15 experimental results – was considered. All experimental data sets are fully rough, and of the DNS data, 30 are considered fully rough flows; all fully rough cases were used for ML training and testing. To the best of our knowledge, this ensemble of roughness geometries is the most extensive used for developing a k_s -prediction method.

In this paper, we first present the governing equations, solution methodologies, simulation parameters and different roughness topographies, and then discuss the post-processed DNS results used to calculate k_s for each surface. Finally, we describe the ML models, their predictions of k_s and their uncertainty.

2. Problem formulation

2.1. Governing equations

The governing equations of incompressible continuity and linear momentum – the Navier–Stokes equations – for a constant-property Newtonian fluid, were solved by DNS. These equations are written in indicial notation as

$$\frac{\partial u_i}{\partial x_i} = 0, \tag{2.1a}$$

$$\frac{\partial u_i}{\partial t} + \frac{\partial u_i u_j}{\partial x_j} = -\frac{\partial P}{\partial x_i} + \nu \frac{\partial^2 u_i}{\partial x_j \partial x_j} + F_i, \tag{2.1b}$$

where $i, j = 1, 2, 3$, x_1, x_2 and x_3 (or x, y, z) are the streamwise, wall-normal and spanwise coordinates, with corresponding velocity components of u_1, u_2 and u_3 (or u, v, w) and P is defined as p/ρ , where p is the pressure and ρ is the fluid density; ν is the kinematic viscosity. An immersed boundary method (Yuan & Piomelli 2014b) was used to enforce the fine-grained roughness boundary conditions on a non-conformal Cartesian grid.

The corresponding body force F_i is added to the right-hand side of the momentum equations to impose a no-slip boundary condition at the fluid–roughness interface. To solve the equations, second-order central differencing was used for spatial discretizations and second-order Adams–Bashforth semi-implicit time advancement was employed. The numerical solver was parallelized using a message passing interface (known as MPI) method (Keating 2004).

A double-averaging decomposition (Raupach & Shaw 1982) was used to resolve turbulent and dispersive components of flow variables in the presence of roughness. In this decomposition, any instantaneous flow variable θ may be decomposed into three components, as

$$\theta(\mathbf{x}, t) = \langle \bar{\theta} \rangle(y) + \theta'(\mathbf{x}, t) + \tilde{\theta}(\mathbf{x}), \quad (2.2)$$

where the time-averaging operator is $\bar{\theta}$ and the intrinsic spatial-averaging operator is $\langle \theta \rangle = 1/A_f \int_{x,z} \theta \, dA$ (and $A_f(y)$ is the area occupied by fluid at an elevation y). The Reynolds and dispersive fluctuating components are then $\theta' = \theta - \bar{\theta}$ and $\tilde{\theta} = \bar{\theta} - \langle \bar{\theta} \rangle$, respectively. Here $\langle \bar{\theta} \rangle$ is called the double-averaged component.

The wall shear stress (including both viscous and pressure drag contributions on a rough wall) was determined by integrating the time-averaged immersed boundary method body force in the x -direction F_1 as

$$\tau_w = \frac{\rho}{L_x L_z} \int_{\mathcal{V}} \overline{F_1}(x, y, z) \, dx \, dy \, dz, \quad (2.3)$$

where \mathcal{V} represents the simulation domain volume below the roughness crest and L_{x_i} is the domain length in the x_i -direction. Readers are referred to Yuan & Piomelli (2014*b,c*) for details of the implementation and validation of the immersed boundary method and the τ_w calculation.

2.2. Surface roughness

In figure 1, surface plots of the 45 roughness geometries used in these simulations are displayed; their statistical properties are given in table 1. Each case name in figure 1 and table 1 begins with the letter C or E, which denotes whether the data is computational or experimental, followed by an identifying index for that particular surface. For computational cases, this index is followed by: a characteristic length scale (as a percentage of δ) used for roughness synthesis; an identifier of whether the surface roughness is regular (reg) or random (rnd); and finally an identifier for one additional surface feature and its position in a series of surfaces with different sizes of that feature. These features were: the streamwise inclination angle I_x in surfaces C01 to C12; the porosity P_o in surfaces C13 to C24; and the streamwise effective slope E_x in surfaces C25 to C30. For the experimental data two indices were assigned to each surface. The first denotes the year in which the data were published and the second is the surface designation in that publication. Thus surfaces with index 16 are from Flack *et al.* (2016), those with index 18 are from Barros, Schultz & Flack (2018) and those with index 19 are from Flack *et al.* (2020). Note that these experimental data were obtained from fully developed channel flows, where the drag was measured through the pressure drop along the channel. Thus their results are expected to be more accurate than those of boundary layer studies where the drag is usually inferred.

Surfaces C01 to C24 were created using ellipsoidal elements (Scotti 2006) of different size, aspect ratio and inclination. For regular roughness, each element had the

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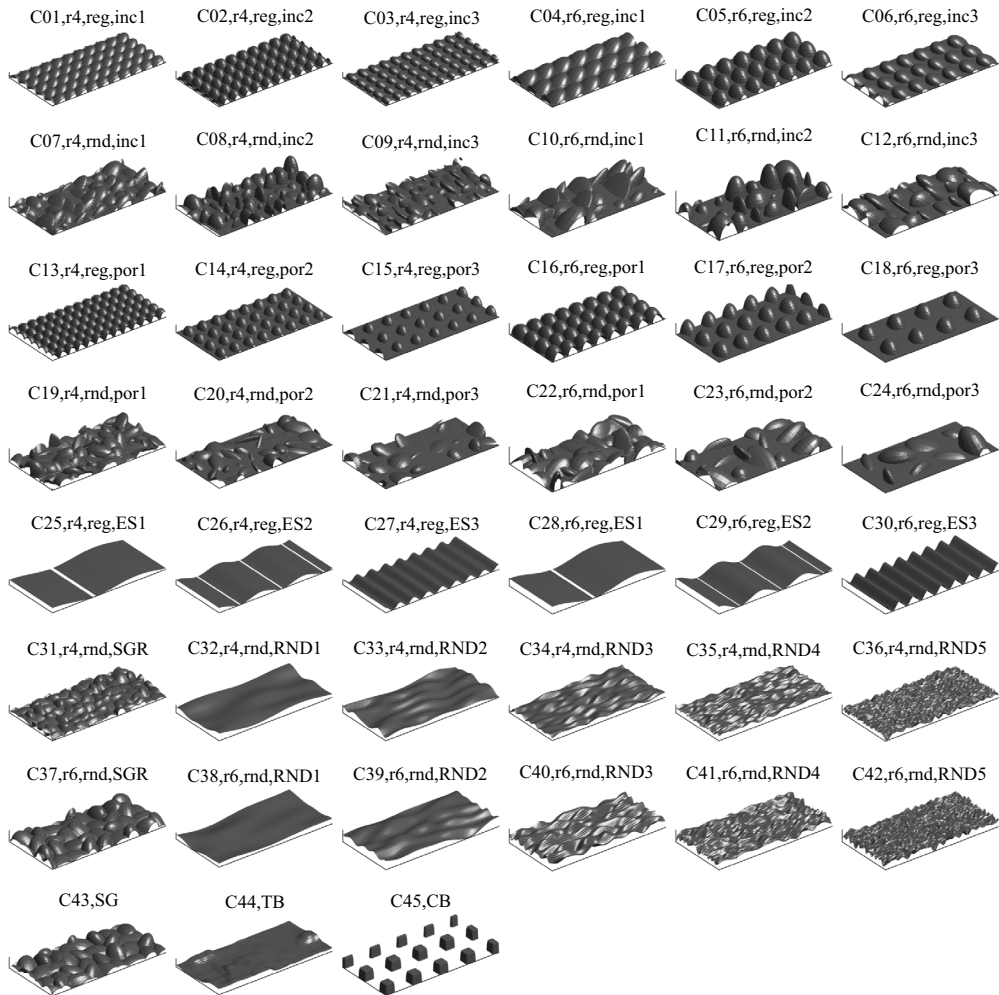


Figure 1. Roughness geometries – each plot is a section of size $\delta \times 0.5\delta$ in the x - z plane. Cases C43 to C45 are from simulations with regular domain sizes (Yuan & Piomelli 2014a; Aghaei Jouybari, Brereton & Yuan 2019).

same orientation and semiaxis lengths, $(\lambda_1, \lambda_2, \lambda_3) = (1.0, 0.7, 0.5)k_c$, where k_c is the peak-to-trough height (also called the crest height). For random roughness, the elements had random orientations and semiaxis lengths (with uniform distributions of the random variables). The average orientation and semiaxis lengths for random roughness were the same as the corresponding regular surface. Surfaces C25 to C30 comprised sinusoidal waves in the x -direction, of the same magnitude but different wavelengths, to generate different values of effective slope E_x . The wavelengths were $3\delta/4$, $3\delta/8$ and $\delta/6$. Surfaces C31 and C37 comprised the random sand-grain roughness of Scotti, which were produced by randomly oriented ellipsoidal elements with fixed semiaxes of $(1.0, 0.7, 0.5)k_c$. Surfaces C32 to C36 and C38 to C42 were generated as the low-order (the first 5, 10, 20, 30 and 50) modes of Fourier transforms of white noise in the streamwise and spanwise directions; they therefore describe random surfaces with large- to small-wavelength roughness. Cases C43, C44 and C45 are DNS results from full-span channel computations

of flow over surfaces of: random sand-grain roughness; the roughness found on a turbine blade (Yuan & Aghaei Jouybari 2018); and arrays of cubes (from the study of Aghaei Jouybari *et al.* (2019)), respectively. Case C46 is a full-span DNS of case C21, generated to validate the minimal-channel approach of the preceding cases. A baseline smooth-wall flow was also simulated using a full-span channel (Yuan & Aghaei Jouybari 2018).

The geometric parameters reported for each surface in table 1 are: roughness peak-to-trough height (also termed crest height) k_c (i.e. distance between the highest and the lowest surface points); mean peak-to-trough height k_t (i.e. the average of peak-to-trough heights obtained from surface tiles of size $\delta \times \delta$, similar to Forooghi *et al.* 2017); mean roughness height k_{avg} ; first-order moment of height fluctuations R_a ; root mean square k_{rms} , skewness S_k and kurtosis K_u of the roughness height fluctuations; surface porosity P_o ; effective slope in the x_i -direction E_{x_i} ; and inclination angle (in radians) in the x_i -direction I_{x_i} , together with the hydrodynamic length scale k_s deduced from the mean velocity field using (1.1).

These geometrical parameters are defined as

$$k_{avg} = \frac{1}{A_t} \int_{x,z} k \, dA, \tag{2.4}$$

$$R_a = \frac{1}{A_t} \int_{x,z} |k - k_{avg}| \, dA, \tag{2.5}$$

$$k_{rms} = \sqrt{\frac{1}{A_t} \int_{x,z} (k - k_{avg})^2 \, dA}, \tag{2.6}$$

$$S_k = \frac{1}{A_t} \int_{x,z} (k - k_{avg})^3 \, dA / k_{rms}^3, \tag{2.7}$$

$$K_u = \frac{1}{A_t} \int_{x,z} (k - k_{avg})^4 \, dA / k_{rms}^4, \tag{2.8}$$

$$E_x = \frac{1}{A_t} \int_{x,z} \left| \frac{\partial k}{\partial x} \right| \, dA, \tag{2.9}$$

$$E_z = \frac{1}{A_t} \int_{x,z} \left| \frac{\partial k}{\partial z} \right| \, dA, \tag{2.10}$$

$$P_o = \frac{1}{A_t k_c} \int_0^{k_c} A_f \, dy, \tag{2.11}$$

$$I_x = \tan^{-1} \left\{ \frac{1}{2} S_k \left(\frac{\partial k}{\partial x} \right) \right\}, \tag{2.12}$$

$$I_z = \tan^{-1} \left\{ \frac{1}{2} S_k \left(\frac{\partial k}{\partial z} \right) \right\}, \tag{2.13}$$

where $k(x, z)$ is the roughness height distribution and $A_f(y)$ and $A_t(y)$ are the fluid and total planar areas at each y location. Here $S_k(\partial k/\partial x_i)$ is the skewness of the $\partial k/\partial x_i$ distribution. In table 1, k_{avg} , k_c , k_{rms} and k_s are then normalized by the first-order moment of height fluctuations R_a and were incorporated in the ML algorithms in this form. All surfaces considered were in the ranges $k_c/\delta \leq 0.17$ and $R_a/\delta \leq 0.04$.

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Case name	k_{avg}	k_c	k_t	k_{rms}	R_a	I_x	I_z	P_o	E_x	E_z	S_k	K_u	k_s
C01,r4,reg,inc1	0.026	0.043	0.043	0.013	0.011	-0.801	-0.089	0.535	0.584	0.510	-0.544	2.177	—
C02,r4,reg,inc2	0.030	0.059	0.059	0.021	0.019	0.012	0.032	0.609	1.029	0.562	-0.265	1.597	—
C03,r4,reg,inc3	0.025	0.043	0.043	0.013	0.011	0.821	-0.078	0.537	0.600	0.485	-0.459	2.052	—
C04,r6,reg,inc1	0.032	0.064	0.064	0.022	0.019	-0.978	0.016	0.597	0.595	0.590	-0.167	1.601	0.064
C05,r6,reg,inc2	0.038	0.088	0.088	0.033	0.030	0.025	0.064	0.654	0.916	0.643	0.109	1.436	0.124
C06,r6,reg,inc3	0.031	0.064	0.064	0.022	0.019	0.955	0.121	0.599	0.588	0.558	-0.087	1.590	0.059
C07,r4,rnd,inc1	0.025	0.086	0.084	0.022	0.019	-0.860	0.033	0.774	0.511	0.559	0.560	2.244	0.136
C08,r4,rnd,inc2	0.027	0.116	0.115	0.030	0.025	-0.007	0.048	0.819	0.861	0.604	0.870	2.627	0.322
C09,r4,rnd,inc3	0.025	0.083	0.081	0.021	0.018	0.829	0.002	0.753	0.517	0.482	0.514	2.292	0.131
C10,r6,rnd,inc1	0.026	0.125	0.120	0.030	0.025	-0.957	-0.019	0.835	0.498	0.578	0.967	2.874	0.269
C11,r6,rnd,inc2	0.033	0.172	0.169	0.044	0.037	0.076	0.138	0.842	0.758	0.543	1.150	3.176	0.536
C12,r6,rnd,inc3	0.032	0.127	0.121	0.032	0.027	0.923	0.032	0.784	0.508	0.471	0.758	2.642	0.272
C13,r4,reg,por1	0.038	0.059	0.059	0.018	0.015	0.024	0.067	0.498	1.043	0.523	-0.820	2.508	—
C14,r4,reg,por2	0.018	0.059	0.059	0.022	0.020	0.021	0.038	0.776	0.613	0.456	0.708	1.840	0.141
C15,r4,reg,por3	0.010	0.059	0.059	0.019	0.014	0.022	0.063	0.877	0.334	0.253	1.646	4.094	0.157
C16,r6,reg,por1	0.051	0.089	0.089	0.030	0.026	0.041	0.149	0.529	1.137	0.534	-0.538	1.873	0.077
C17,r6,reg,por2	0.022	0.089	0.089	0.031	0.027	0.041	0.080	0.801	0.537	0.403	0.982	2.308	0.260
C18,r6,reg,por3	0.013	0.089	0.089	0.026	0.020	0.057	0.126	0.886	0.307	0.230	1.849	4.839	0.247
C19,r4,rnd,por1	0.027	0.112	0.108	0.021	0.017	0.025	-0.107	0.806	0.487	0.486	0.732	3.422	0.158
C20,r4,rnd,por2	0.013	0.095	0.087	0.017	0.014	0.032	-0.646	0.896	0.311	0.323	1.343	4.126	0.106
C21,r4,rnd,por3	0.009	0.098	0.094	0.016	0.012	0.321	-0.741	0.929	0.219	0.233	2.168	7.728	0.103
C22,r6,rnd,por1	0.035	0.139	0.139	0.029	0.024	-0.070	-0.245	0.791	0.456	0.499	0.591	2.830	0.277
C23,r6,rnd,por2	0.017	0.123	0.111	0.025	0.020	-0.672	-0.841	0.885	0.305	0.325	1.467	4.347	0.175
C24,r6,rnd,por3	0.014	0.152	0.145	0.027	0.019	0.189	-0.056	0.926	0.254	0.257	2.371	8.740	0.260
C25,r4,reg,ES1	0.020	0.040	0.040	0.014	0.013	0.046	0.006	0.510	0.106	0.009	-0.032	1.503	—
C26,r4,reg,ES2	0.021	0.040	0.040	0.014	0.013	0.039	-0.001	0.510	0.212	0.020	-0.071	1.505	0.065
C27,r4,reg,ES3	0.023	0.040	0.040	0.014	0.012	0.006	-0.023	0.510	0.609	0.032	-0.214	1.544	—
C28,r6,reg,ES1	0.030	0.059	0.059	0.021	0.019	0.044	0.018	0.504	0.158	0.015	-0.031	1.499	0.071
C29,r6,reg,ES2	0.031	0.059	0.059	0.021	0.019	0.028	-0.069	0.504	0.316	0.022	-0.071	1.503	0.112
C30,r6,reg,ES3	0.034	0.059	0.059	0.020	0.018	0.015	-0.069	0.505	0.917	0.048	-0.203	1.543	0.064
C31,r4,rnd,SGR	0.025	0.059	0.059	0.011	0.009	0.104	-0.039	0.648	0.370	0.398	0.378	2.784	0.049
C32,r4,rnd,RND1	0.040	0.075	0.072	0.013	0.010	0.117	0.108	0.479	0.068	0.169	-0.069	2.991	—
C33,r4,rnd,RND2	0.041	0.088	0.084	0.013	0.011	0.109	0.078	0.553	0.117	0.308	0.004	2.763	—
C34,r4,rnd,RND3	0.042	0.080	0.071	0.010	0.008	0.070	0.051	0.508	0.175	0.458	-0.002	3.031	—
C35,r4,rnd,RND4	0.043	0.077	0.066	0.008	0.007	0.039	0.042	0.488	0.218	0.558	0.013	2.941	—
C36,r4,rnd,RND5	0.045	0.084	0.067	0.009	0.007	0.035	0.037	0.535	0.378	0.841	0.075	3.018	—
C37,r6,rnd,SGR	0.037	0.088	0.088	0.018	0.015	0.312	0.180	0.640	0.428	0.463	0.323	2.686	0.109
C38,r6,rnd,RND1	0.060	0.106	0.091	0.016	0.012	0.045	0.028	0.444	0.077	0.183	-0.220	3.258	—
C39,r6,rnd,RND2	0.061	0.098	0.095	0.012	0.009	0.111	0.057	0.400	0.108	0.285	-0.020	3.267	—
C40,r6,rnd,RND3	0.064	0.121	0.112	0.016	0.013	0.061	0.022	0.512	0.280	0.760	0.037	2.977	0.050
C41,r6,rnd,RND4	0.065	0.130	0.130	0.015	0.012	0.045	0.037	0.546	0.374	0.989	0.028	3.036	—
C42,r6,rnd,RND5	0.068	0.118	0.116	0.013	0.010	0.037	0.025	0.503	0.547	1.204	0.052	2.933	—
C43,SG	0.036	0.089	0.087	0.017	0.014	0.288	0.156	0.649	0.425	0.441	0.476	2.970	0.093
C44,TB	0.055	0.125	0.088	0.018	0.014	0.007	-0.006	0.569	0.097	0.081	0.200	3.493	0.024
C45,CB	0.010	0.070	0.070	0.023	0.016	0.420	0.508	0.878	0.249	0.247	2.101	5.569	0.150
C46,r4,rnd,por3,FS	0.009	0.098	0.094	0.016	0.012	0.321	-0.715	0.929	0.219	0.234	2.168	7.728	0.104
E01,16,2	0.138	0.261	0.254	0.020	0.016	-0.005	0.011	0.472	0.720	0.835	-0.711	3.843	0.052
E02,16,3	0.143	0.252	0.252	0.021	0.016	-0.021	0.010	0.432	0.740	0.868	-0.338	3.159	0.050
E03,16,7	0.133	0.365	0.254	0.019	0.014	-0.038	0.000	0.638	0.618	0.705	-1.169	5.292	0.058
E04,16,8	0.126	0.298	0.227	0.017	0.013	-0.034	0.009	0.579	0.587	0.682	-1.445	5.421	0.056
E05,16,9	0.112	0.308	0.167	0.018	0.014	-0.031	0.015	0.637	0.636	0.753	-0.738	3.714	0.043
E06,16,15	0.081	0.191	0.191	0.013	0.010	-0.027	0.003	0.578	0.621	0.713	-0.687	3.854	0.035
E07,18,1	0.121	0.241	0.227	0.026	0.021	-0.013	-0.183	0.500	0.181	0.188	0.107	2.941	0.053
E08,18,2	0.143	0.276	0.255	0.032	0.025	-0.019	0.194	0.483	0.162	0.164	0.093	2.967	0.034
E09,19,1	0.204	0.398	0.344	0.046	0.036	0.042	-0.096	0.487	0.227	0.230	-0.080	2.989	0.065
E10,19,2	0.389	0.763	0.689	0.088	0.070	0.046	0.002	0.492	0.447	0.452	-0.065	2.925	0.200
E11,19,3	0.477	0.730	0.679	0.088	0.070	-0.029	-0.245	0.348	0.434	0.432	-0.660	3.274	0.160
E12,19,4	0.459	0.751	0.710	0.089	0.071	-0.052	0.036	0.391	0.455	0.459	-0.351	3.041	0.180
E13,19,5	0.292	0.732	0.650	0.090	0.072	-0.058	-0.004	0.602	0.445	0.452	0.346	3.051	0.245
E14,19,6	0.202	0.711	0.604	0.087	0.069	0.004	-0.010	0.716	0.391	0.400	0.812	3.559	0.435
E15,19,7	0.522	0.967	0.894	0.114	0.092	-0.050	-0.235	0.462	0.557	0.562	-0.066	2.794	0.230

Table 1. Statistical parameters of roughness topography and the equivalent sand-grain height k_s for each roughness geometry. Here R_a , k_{avg} , k_c , k_t , k_{rms} and k_s values from DNS are normalized by the channel half-height δ , while corresponding experimental values are given in mm; k_s is not listed for cases thought to be transitionally rough.

2.3. Simulation parameters

Direct numerical simulation was used to calculate the velocity and pressure fields in turbulent open-channel flows over 45 different rough surfaces and one smooth one, at a constant frictional Reynolds number $Re_\tau = u_\tau \delta / \nu = 1000$, where u_τ is the friction velocity and δ is the channel half-height. In these simulations, the domain sizes were $(L_x, L_y, L_z) = (3, 1, 1)\delta$. The origin of the y axis was the elevation of the lowest trough for each rough surface. The number of grid points was $(n_x, n_y, n_z) = (400, 300, 160)$. A uniform mesh was used in the x - and z -directions, yielding grid sizes of $\Delta x^+ = 7.5$ and $\Delta z^+ = 6.3$, where $+$ denotes normalization in wall units. For all cases, the mesh was stretched in the y -direction with a hyperbolic tangent function, with the third grid point from the origin at $y^+ < 1$. For the rough-wall cases, at the roughness crest, $\Delta y/k_c \leq 0.017$, with this ratio taking its highest value for case C11. The maximum grid size was $\Delta y_{max}^+ = 9.5$ at the channel centreline, where the Kolmogorov length scale $\eta^+ \approx 6$. Moin & Mahesh (1998) have proposed that one requirement for obtaining reliable first- and second-order flow statistics is that the grid resolution must be fine enough to capture accurately most of the dissipation, while Moser & Moin (1987) noted that most of the dissipation in curved channel flow occurs at scales greater than 15η (based on average dissipation). It follows that for DNS computations of these kinds of flow statistics in channel and boundary-layer flows, $\Delta x/\eta$ and $\Delta z/\eta$ are typically chosen between 7 to 15 and 4 to 8, respectively (see, for example Kim, Moin & Moser 1987; Spalart 1988; Yuan & Piomelli 2014c). The grid sizes in this study were chosen accordingly and were $\Delta x/\eta < 7.5$, $\Delta y/\eta < 4.0$ and $\Delta z/\eta < 6.5$.

Periodic boundary conditions were imposed in the streamwise and spanwise directions, with no-slip and symmetry boundary conditions at the bottom and top boundaries, respectively. After each simulation had reached statistical stationarity, data were collected for ensemble averaging over 10 large-eddy turn-over times (δ/u_τ). In these simulations, the time step $\tau^+ \leq 0.04$ and so was significantly smaller than the largest acceptable one of $\tau^+ \approx 0.2$ recommended by Choi & Moin (1994) for DNS.

The surface Taylor microscales $\lambda_{T,x}$ and $\lambda_{T,z}$, in the x - and z -directions, were used to evaluate the adequacy of the grid resolution for resolving details of flow in the roughness sublayer, following Yuan & Piomelli (2014b). These geometric microscales were obtained by fitting a parabola to the two-point autocorrelation of the surface height fluctuation in the respective direction. They represent the size of an equivalent ‘roughness element’ in the context of random multiscale roughness. The streamwise and spanwise values of λ_T , rescaled by u_τ/ν as λ_T^+ , and the respective grid sizes are given in table 2 (part I). For each case, λ_{T,x_i}^+ is of order 10 to 10^2 , indicating that the average size of the roughness element is large in viscous units. On average, roughness elements were well resolved by the grid, with typically 4 to 12 grid points per λ_{T,x_i} microscale in each direction. For reference purposes, Yuan & Piomelli (2014a) reported a resolution of $\lambda_{T,x}/\Delta x \approx 4$ in their large-eddy simulations of channel flow over surfaces with sand-grain roughness. The cases in table 2 for which λ_T was not well resolved in at least one direction ($\lambda_{T,x}/\Delta x < 3$ or $\lambda_{T,z}/\Delta z < 3$) may also not have been fully rough flows (as discussed in the following section), and so were not included in the ensemble of flows for ML training and testing.

In rough-wall flows, the pressure drag is caused primarily by the local flow structures and separation in the vicinity of individual roughness protuberances, which are predominately near-wall phenomena. To carry out the 46 separate DNS simulations for determining k_s efficiently, with sufficient near-wall resolution, a small-span channel simulation approach was employed. The concept of minimal-span simulation was introduced by Jimenez & Moin (1991). Chung *et al.* (2015) and MacDonald *et al.* (2017) carried out analyses of the performance of DNS over small spanwise domains for full and

Case name	Part I				Part II	
	$\lambda_{T,x}^+$	$\lambda_{T,x}/\Delta x$	$\lambda_{T,z}^+$	$\lambda_{T,z}/\Delta z$	d/δ	\hat{k}_s^+
C01,r4,reg,inc1	19.7	2.6	21.1	3.4	0.032	19.4
C02,r4,reg,inc2	20.4	2.7	33.1	5.3	0.046	49.7
C03,r4,reg,inc3	19.8	2.6	22.9	3.7	0.033	31.0
C04,r6,reg,inc1	27.7	3.7	28.4	4.5	0.038	64.4
C05,r6,reg,inc2	31.6	4.2	39.1	6.2	0.057	124.4
C06,r6,reg,inc3	29.9	4.0	30.0	4.8	0.045	58.9
C07,r4,rnd,inc1	33.8	4.5	26.7	4.3	0.036	136.2
C08,r4,rnd,inc2	26.1	3.5	32.7	5.2	0.052	322.3
C09,r4,rnd,inc3	35.5	4.7	30.1	4.8	0.039	131.1
C10,r6,rnd,inc1	38.2	5.1	29.7	4.8	0.042	268.9
C11,r6,rnd,inc2	38.1	5.1	47.0	7.5	0.070	536.4
C12,r6,rnd,inc3	47.9	6.4	40.2	6.4	0.053	271.7
C13,r4,reg,por1	17.8	2.4	32.7	5.2	0.047	41.4
C14,r4,reg,por2	27.5	3.7	34.2	5.5	0.032	140.6
C15,r4,reg,por3	31.5	4.2	39.4	6.3	0.028	157.1
C16,r6,reg,por1	25.6	3.4	46.1	7.4	0.066	76.7
C17,r6,reg,por2	40.1	5.3	47.8	7.6	0.044	259.8
C18,r6,reg,por3	44.4	5.9	54.8	8.8	0.039	246.5
C19,r4,rnd,por1	32.7	4.4	31.1	5.0	0.042	158.2
C20,r4,rnd,por2	35.6	4.7	31.3	5.0	0.026	105.7
C21,r4,rnd,por3	37.4	5.0	34.2	5.5	0.027	102.7
C22,r6,rnd,por1	44.6	5.9	35.3	5.6	0.053	276.8
C23,r6,rnd,por2	47.1	6.3	39.7	6.4	0.038	175.1
C24,r6,rnd,por3	47.1	6.3	44.4	7.1	0.045	260.3
C25,r4,reg,ES1	89.0	11.9	—	—	0.024	25.6
C26,r4,reg,ES2	66.5	8.9	—	—	0.026	65.3
C27,r4,reg,ES3	27.1	3.6	—	—	0.035	45.5
C28,r6,reg,ES1	90.6	12.1	—	—	0.033	71.2
C29,r6,reg,ES2	66.8	8.9	—	—	0.040	112.0
C30,r6,reg,ES3	27.2	3.6	—	—	0.054	64.0
C31,r4,rnd,SGR	27.8	3.7	25.0	4.0	0.032	48.7
C32,r4,rnd,RND1	131.2	17.5	54.1	8.7	0.041	8.4
C33,r4,rnd,RND2	96.3	12.8	42.1	6.7	0.043	17.6
C34,r4,rnd,RND3	56.4	7.5	22.4	3.6	0.045	22.5
C35,r4,rnd,RND4	39.5	5.3	15.8	2.5	0.046	18.3
C36,r4,rnd,RND5	25.1	3.3	11.4	1.8	0.051	23.4
C37,r6,rnd,SGR	36.5	4.9	31.9	5.1	0.046	108.8
C38,r6,rnd,RND1	88.5	11.8	72.6	11.6	0.060	12.0
C39,r6,rnd,RND2	93.8	12.5	35.7	5.7	0.062	17.1
C40,r6,rnd,RND3	57.0	7.6	22.8	3.6	0.070	50.4
C41,r6,rnd,RND4	40.5	5.4	15.6	2.5	0.073	48.7
C42,r6,rnd,RND5	24.5	3.3	11.3	1.8	0.076	43.8
C43,SG	35.2	6.0	33.5	5.7	0.044	93.0
C44,TB	132.1	10.4	168.5	13.2	0.058	24.1
C45,CB	25.7	4.5	25.5	4.4	0.039	149.9
C46,r4,rnd,por3,FS	37.6	5.0	34.6	5.5	0.027	104.2

Table 2. Part I: streamwise and spanwise values of the surface Taylor microscale λ_T . Part II: flow-related parameters obtained from DNS. The flow is assumed fully rough if $\hat{k}_s^+ \gtrsim 50$, in which case k_s is equal to \hat{k}_s .

open channel flows on rough and smooth walls and showed that minimal-span simulations captured the essential near-wall dynamics and yielded accurate computations of wall friction, and of mean velocities and Reynolds stresses as far from the wall as $y \approx 0.3\delta$, when the following constraints were met:

$$L_x \geq \max(1000\delta_v, 3L_z, \lambda_{r,x}), \tag{2.14a}$$

$$L_y \geq k_c/0.15, \tag{2.14b}$$

$$L_z \geq \max(100\delta_v, k_c/0.4, \lambda_{r,z}), \tag{2.14c}$$

where $\delta_v = \nu/u_\tau$ and λ_{r,x_i} is the characteristic roughness wavelength in the x_i -direction. Alternatively, the surface Taylor microscale may be used as the length scale in this constraint. Conditions (2.14a,c) were satisfied by choosing domain sizes L_x^+ and L_z^+ of 3000 and 1000, respectively, while condition (2.14b) was met for all cases except C11, which fell below the $L_y \geq k_c/0.15$ constraint by approximately 10% – C11 is a case with random geometry; protuberances beyond 0.15δ exist but are rare.

The criteria of (2.14) were developed originally for simulations of flow over surfaces with uniformly distributed roughness elements. In this study, the random roughness geometries used require an additional criterion on the sufficiency of the domain size: the area $L_x L_z$ should be large enough to achieve statistical convergence of surface parameters, such as k_{rms} and E_{x_i} , and of the flow parameter k_s . To check the adequacy of the chosen domain size, an additional simulation was carried out of case C21, the surface comprising the largest dominant spatial wavelength (and consequently the most limited sampling of random geometrical components with this wavelength) and a long-tailed height-fluctuation probability density function (p.d.f.) with a kurtosis of around 8. In this validation simulation, denoted case C46, the domain sizes were doubled in x and z , by duplicating C21 in these directions. The double-averaged velocity profiles $U^+ = \langle \bar{u} \rangle^+(y^+)$ for cases C21 and C46 are in a very good agreement over the log-linear region, as shown in figure 2. Each surface statistic differs by no more than 3%, with the greatest discrepancy found in I_z , while the equivalent sand-grain roughness height k_s is almost equal in the two cases. The chosen domain size was therefore considered sufficient for accuracy and convergence of statistics describing flow over the random roughness geometries of this study.

3. Results

3.1. Post-processed results

In figure 2, the streamwise double-averaged velocity profiles computed in these simulations are shown. The profiles in the logarithmic region are described for the smooth-wall case and the fully rough rough-wall cases as

$$\langle \bar{u} \rangle^+ = \frac{1}{\kappa} \ln(y^+) + 5.0 \quad \text{and} \tag{3.1a}$$

$$\langle \bar{u} \rangle^+ = \frac{1}{\kappa} \ln\left(\frac{y-d}{k_s}\right) + 8.5, \tag{3.1b}$$

respectively, where d is the zero-plane displacement, obtained as the location of the centroid of the wall-normal profile of the averaged drag force (Jackson 1981). The shift in the y coordinate by d accounts for the flow blockage by surface roughness elements, and the values of d are given in table 2 (part II).

Equivalent sand-grain height prediction of rough walls

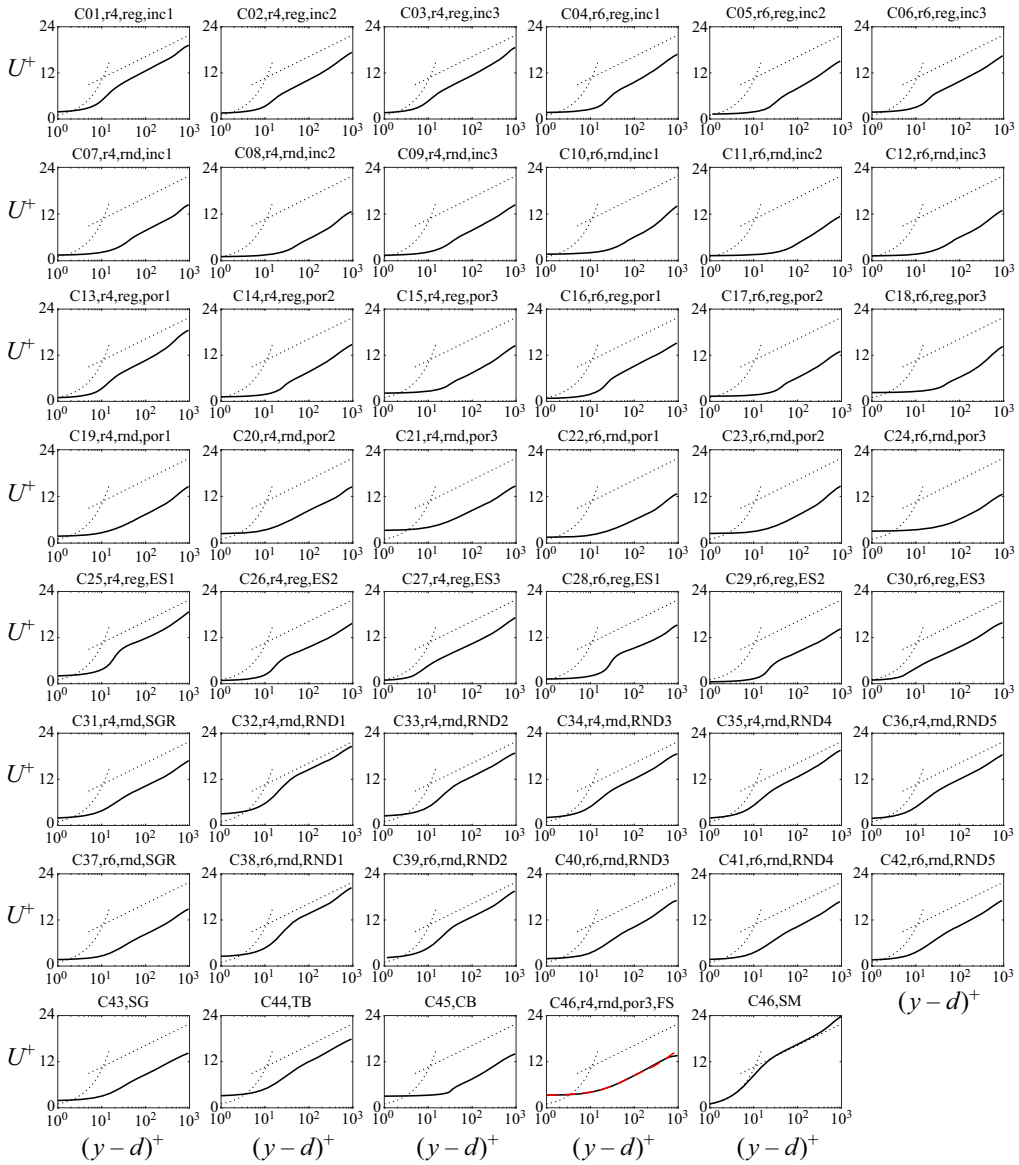


Figure 2. Profiles of streamwise double-averaged velocity plotted against a zero-plane-displacement shifted logarithmic y abscissa. The dashed lines are $u^+ = y^+$ and $u^+ = 2.5 \ln(y-d)^+ + 5.0$. The red dot-dash line in plot C46 is that of C21.

To determine whether a particular flow was within the fully rough regime, (3.1b) was applied to the computed logarithmic velocity profile to yield a test value of k_s , denoted as \hat{k}_s in table 2 (part II). With \hat{k}_s determined for all cases, those with \hat{k}_s^+ greater than a threshold value of 50 were deemed to be in the fully rough regime (30 surfaces), in which case k_s was set to equal \hat{k}_s . Those below the threshold were possibly transitionally rough (15 surfaces) and so were not included in ML predictions in this study. The threshold value of k_s^+ – the lower end of the fully rough regime – has been observed to vary significantly for different types of roughness and is typically between 20 and 80. For example, the threshold values for surfaces C43 and C44 are roughly 80 and 20 (Yuan & Piomelli 2014a) and 50 for surface C45 (Bandyopadhyay 1987).

The threshold value of k_s^+ which signifies the beginning of the fully rough regime was not determined more precisely because of the cost of carrying out, for each surface, simulations at successively higher values of k_s^+ until k_s/R_a became invariant with the Reynolds number. In the GPR prediction, potential uncertainties in k_s which might arise through treating all flows with $k_s^+ > 50$ as fully rough, and other sources of possible error, were compensated for by incorporating an assumed 10 % noise level in the learning stage of the prediction of k_s , as discussed in § 3.2. The values of $k_s^+ = 50$ as the threshold for fully rough flows and the assumed noise level were chosen as part of a trade-off to maximize the number of usable data, to avoid overfitting, while acknowledging possible uncertainties in the modelling data.

In figure 3, pair plots of the different topographic roughness parameters are shown as scatter plots (lower left), joint p.d.f.s (upper right) and distribution p.d.f.s (diagonal). Pair scatter plots for the true (DNS and experimental) value of k_s and other roughness parameters are along the bottom row of this figure. It can be seen that, for the roughness cases chosen, there is some correlation between kurtosis and r.m.s. roughness (column 1, row 6), kurtosis and skewness (column 5, row 6) and skewness and porosity (column 2, row 5). The relationship between others appears to be more random. From the graphs in the bottom row, it can be seen that k_s/R_a scales on porosity to some power, albeit with some scatter (column 2, row 7). It also appears that k_s/R_a might decrease with skewness for surfaces with $S_k < 0$ and increase with skewness in cases with $S_k > 0$ (column 5, row 7). Surfaces with positive skewness yielded higher values of k_s compared with those with negative skewness, consistent with the observation of Flack *et al.* (2020). Beyond these observations, there does not appear to be a clear linear correlation between k_s and any individual roughness parameter, which makes the search for a functional dependence of k_s on these parameters a problem well suited to ML. The measures of inclination, I_x and I_z , showed no clear correlation with other variables or with k_s/R_a .

3.2. ML predictions of the equivalent sand-grain height

The ML techniques of DNN and GPR were employed to predict k_s from the data sets described in the previous section. The objectives of this exercise were to generate and collect data, and make qualitative comparisons between ML predictions and those from conventional correlations, rather than evaluating and comparing the performance of various ML procedures *per se*. The DNN and GPR approaches were used because our experience was that they predicted k_s with high accuracy, notwithstanding their simplicity. Other approaches such as the support vector machine technique were considered initially, but their preliminary predictions were not as accurate as those found using DNN and GPR approaches.

The main characteristics of DNN and GPR methods are described below.

- (i) The inputs for both techniques were 17 roughness geometrical parameters, eight of which were the primary variables k_{rms}/R_a , I_x , $|I_z|$, P_o , E_x , E_z , S_k and K_u (defined in (2.4) to (2.13)). The other nine were products of the primary variables, which were added to improve the efficiency of each learning stage. They were $p_1 = E_x E_z$, $p_2 = E_x S_k$, $p_3 = E_x K_u$, $p_4 = E_z S_k$, $p_5 = E_z K_u$, $p_6 = S_k K_u$, $p_7 = E_x^2$, $p_8 = E_z^2$ and $p_9 = S_k^2$. These particular products were chosen because of their perceived importance for certain types of roughness.
- (ii) The database consisted of 45 different sets: 30 DNS of turbulent channel flows over different surfaces at $Re_\tau = 1000$, and 15 experimental data sets at higher Reynolds numbers, with all data sets in the fully rough turbulent-flow regime.

Equivalent sand-grain height prediction of rough walls

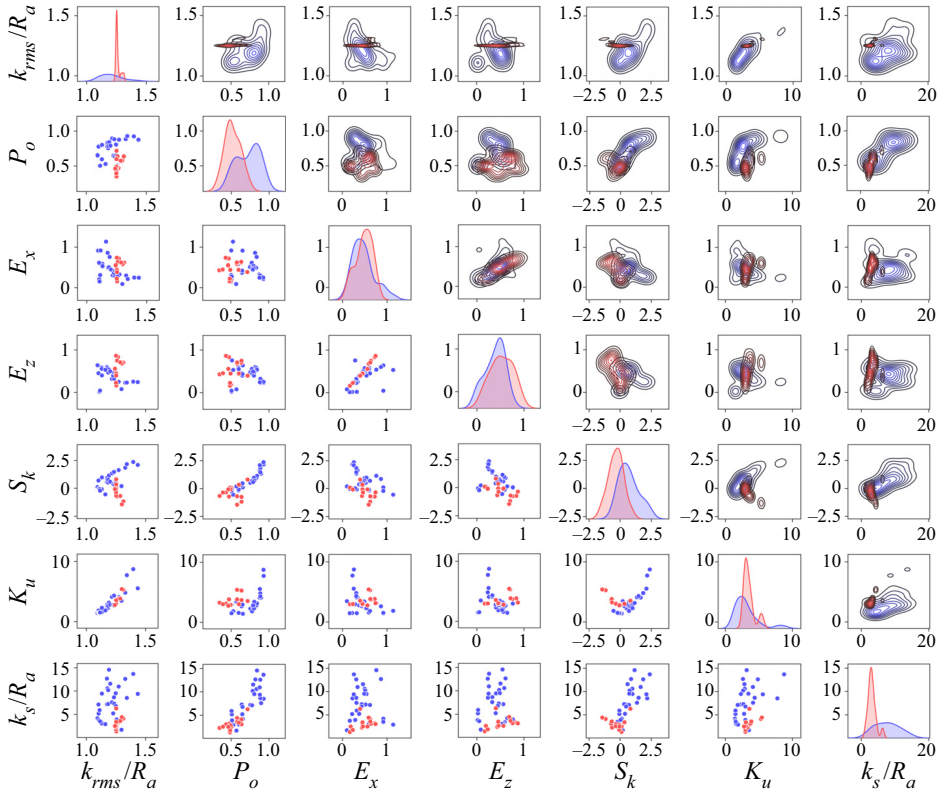


Figure 3. Pair plots of geometrical parameters and k_s , with k_s plots in the bottom row and the first column, DNS data (blue), experimental data (red).

- (iii) The DNN architecture was a multilayer perceptron, with three hidden layers (with 18, 7 and 7 neurons, respectively). The activation functions at all nodes were of the rectified linear unit kind, and kernel regularization was used to avoid overfitting. The network had 521 trainable weights in total. The preset parameters to the algorithm were optimized based on available data, through a hyperparameter tuning process. Specifically, 270 configurations were first generated with different lengths (representing the number of layers) and widths (representing the number of neurons). For each configuration, the DNN compiler was performed 1000 times with random selections of training (70 % of total) and testing (30 % of total) datasets to identify the best performance of the configuration. The configuration that yielded the best results was considered as the optimal one, the results of which are presented here. The cost of data fitting for one iteration (out of 1000) for each DNN configuration was approximately one second. In total, it took approximately 75 hours to obtain the optimal DNN network. This architecture was found to provide suitable accuracy in modelling without overfitting, for this particular multivariate labelled regression problem.
- (iv) The GPR procedure used rational quadratic kernels to represent k_s as a superposition of scaled Gaussian functions of the independent variables of the modelling problem. Similar to the DNN method, the training and testing data were chosen randomly, with respective ratios of 70 % and 30 % of the total data points. The preset parameters

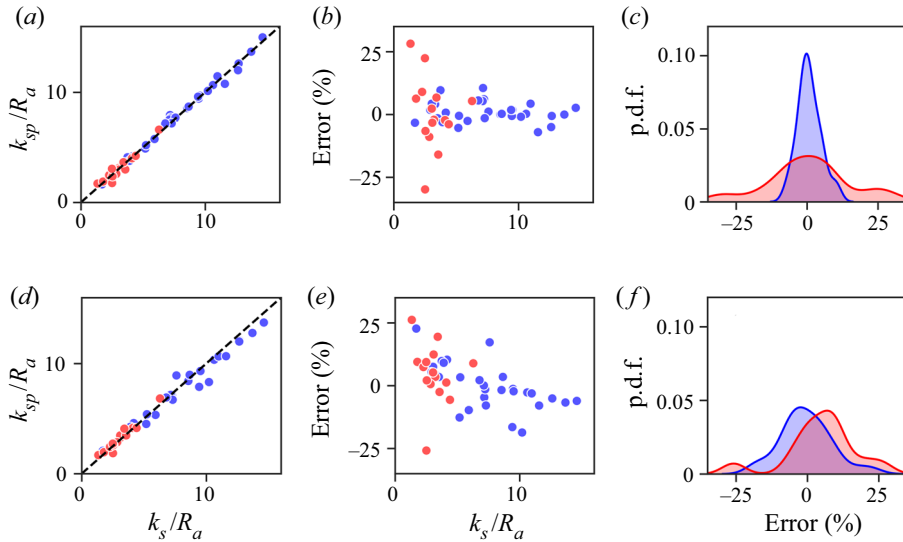


Figure 4. (a,d) Scatter plot of true k_s and predicted k_s , (b,e) scatter plot of true k_s and relative error, (c,f) p.d.f.s of relative error, for (a–c) DNN and (d–f) GPR predictions, with DNS data (blue), experimental data (red).

(e.g. kernel type, number of iterations, etc.) were also tuned with the available data by running the GPR compiler approximately 8000 times. It took approximately 35 hours to obtain the optimal fit. The GPR method has the capability of incorporating uncertainty or noise in the determination of model parameters in the learning stages. Such noise might arise through: numerical and discretization errors; uncertainty in the form and model coefficients of equation (1.1); the applicability and fitting range of equation (1.1) (which was deduced from high Reynolds number experiments) to simulations at much lower Reynolds numbers; and the possibility that some of the training data may have been from simulations in which the flow was not quite fully rough. A noise level of 10% in k_s/R_a values was chosen as an upper estimate of the likely uncertainty from these sources. Noise levels of 5% and 15% were also tested, but little sensitivity of the k_s prediction was found to the assumed noise level within the tested range.

The values of k_s predicted from the surface topography parameters, henceforth called k_{sp} , are compared with the actual k_s values in figure 4, for the DNN and GPR methods, respectively. Scatter plots of k_{sp} and the true value of k_s in figures 4(a) and 4(d) reveal a tight clustering of data along the $y = x$ diagonal, with only a few outlying points. This very high degree of correlation between k_{sp} and k_s implies that both techniques have been applied with equal success to this prediction problem. The error range, figures 4(b) and 4(e), is less than $\pm 30\%$ (L_∞ norm) and the average error (L_1 norm) is less than 8%, for both techniques.

The consistency between both the k_s predictions and error bands for two quite different ML techniques suggests that they are both well-suited to this kind of problem, and possibly close to an optimum for this class of ML approach.

The error values as percentages, for the DNN and GPR methods, are given in table 3, together with the error in the empirical relation

$$k_s = 2.91k_{rms}(2 + S_k)^{-0.284}, \tag{3.2}$$

Case name	err_{DNN}	err_{GPR}	err_{B1}	err_{B2}	err_{B3}	err_{B4}
C04,r6,reg,inc1	4.0	4.1	-16.7	-40.9	8.6	-63.9
C05,r6,reg,inc2	0.7	10.3	-38.3	-49.5	2.0	-71.7
C06,r6,reg,inc3	4.2	7.5	-10.4	-33.6	24.5	-59.8
C07,r4,rnd,inc1	10.5	-4.7	-63.5	-63.6	10.0	-73.5
C08,r4,rnd,inc2	-0.6	-4.8	-80.1	-77.6	-4.1	-81.7
C09,r4,rnd,inc3	6.0	-1.5	-63.4	-64.2	8.3	-73.4
C10,r6,rnd,inc1	0.2	-2.7	-76.3	-72.5	11.8	-77.8
C11,r6,rnd,inc2	2.6	-6.1	-82.9	-78.9	4.1	-82.2
C12,r6,rnd,inc3	-1.0	-18.7	-74.7	-72.7	-2.3	-78.7
C14,r4,reg,por2	5.3	0.0	-66.2	-64.2	-8.7	-80.3
C15,r4,reg,por3	4.2	-3.1	-76.7	-66.5	29.0	-78.8
C16,r6,reg,por1	1.8	5.4	3.5	-41.7	21.5	-59.4
C17,r6,reg,por2	-0.6	-1.3	-74.5	-70.3	-10.9	-82.7
C18,r6,reg,por3	-5.1	-5.1	-79.1	-68.4	35.5	-78.8
C19,r4,rnd,por1	1.8	-2.3	-71.4	-69.4	44.9	-67.8
C20,r4,rnd,por2	1.1	17.2	-67.0	-56.6	82.1	-66.3
C21,r4,rnd,por3	0.0	-1.8	-69.6	-50.0	254.1	-46.2
C22,r6,rnd,por1	-7.1	-7.9	-77.0	-76.7	-10.6	-78.4
C23,r6,rnd,por2	0.2	3.4	-70.9	-60.4	80.8	-67.9
C24,r6,rnd,por3	-0.1	-6.7	-80.5	-66.3	136.7	-66.5
C26,r4,reg,ES2	-5.4	-12.7	-48.6	-61.6	-57.6	-83.8
C28,r6,reg,ES1	9.6	9.8	-29.2	-45.9	-51.9	-81.2
C29,r6,reg,ES2	-2.6	-9.8	-54.7	-66.2	-53.2	-83.2
C30,r6,reg,ES3	-1.5	3.4	-21.8	-45.7	8.1	-65.7
C31,r4,rnd,SGR	-0.6	3.3	-46.7	-50.7	65.1	-53.8
C37,r6,rnd,SGR	-1.5	-7.9	-61.3	-65.0	11.9	-68.6
C40,r6,rnd,RND3	-3.1	9.1	-23.6	-39.6	98.3	-30.8
C43,SG	5.5	2.1	-58.6	-60.1	46.3	-62.0
C44,TB	-3.3	22.7	77.6	51.9	31.5	-51.6
C45,CB	1.8	-16.5	-70.4	-52.0	79.3	-72.8
E01,16,2	-2.1	3.5	6.2	-47.5	370.2	63.0
E02,16,3	2.3	5.2	3.3	-33.7	429.4	79.5
E03,16,7	-2.3	1.2	-2.2	-69.1	368.1	38.6
E04,16,8	-3.9	-5.7	1.3	-78.8	412.4	27.6
E05,16,9	-3.3	12.4	10.9	-46.3	262.1	27.3
E06,16,15	-16.0	-2.5	-3.0	-51.1	405.4	79.9
E07,18,1	-29.8	-25.8	17.3	-4.0	208.3	11.2
E08,18,2	28.1	26.1	120.7	79.4	388.8	80.0
E09,19,1	6.2	9.4	69.2	25.9	312.5	56.9
E10,19,2	-8.9	0.6	5.8	-20.7	258.9	20.6
E11,19,3	8.9	7.4	47.4	-24.1	247.4	32.2
E12,19,4	-6.6	2.1	24.1	-21.0	258.4	32.2
E13,19,5	6.7	19.4	-16.6	-23.8	287.2	6.6
E14,19,6	5.3	8.9	-56.8	-52.5	177.2	-38.2
E15,19,7	22.3	9.4	19.8	-10.2	342.6	43.0
L_1	5.4	7.8	47.6	52.8	133.8	60.6
L_∞	29.8	26.1	120.7	79.4	429.4	83.8

Table 3. Errors in k_s prediction by DNN and GPR compared with errors of the empirical correlations: err_{B1} (3.2), err_{B2} (3.4), err_{B3} (3.3) and err_{B4} (3.5). The four largest errors (in magnitude) for each column are coloured in red. The errors are percentages.

proposed by Flack *et al.* (2016) and

$$k_s = 1.07k_t(1 - e^{-3.5E_x})(0.67S_k^2 + 0.93S_k + 1.3), \quad (3.3)$$

given by Forooghi *et al.* (2017), as well as their respective recalibrated correlations

$$k_s = 1.11k_{rms}(2 + S_k)^{0.74}, \quad (3.4)$$

$$k_s = 0.04k_t(1 - e^{-5.50E_x})(S_k^2 + 2.57S_k + 9.82), \quad (3.5)$$

when extended to all cases in the current database. It is interesting to note that the form of equation (3.2) was chosen for surfaces generated by grit blasting – closely packed, random, three-dimensional (3-D) roughness with a wide range of scales (E01–E06), while many of the simulated surfaces are two-dimensional (2-D), some are characterized by discrete elements of similar sizes, while others are sparse or wavy (characterized by low slopes). Equation (3.3), on the other hand, includes a slope parameter and was calibrated for numerically generated surfaces consisting elements of random sizes and a prescribed shape.

For most cases, the errors from the DNN and GPR methods were of the same order of magnitude and much smaller than the error in using (3.2) or (3.3). In the DNN and GPR predictions of simulation cases, the greatest errors (approximately 25%–30%) arose in cases E07 and E08. The surfaces associated with these cases are characterized by fractal features (with spectral slopes of -0.5 and -1.0 , respectively (Barros *et al.* 2018)). The size of the errors for these cases might be attributed to the small number of surfaces with this feature used in the training set (as opposed to the many surfaces that are mostly characterized by single-scale elements). A close examination of the prediction errors for the DNS cases showed a subtle trend between relatively high errors and low roughness solidity (or low E_s and insignificant wake sheltering), in, for example, cases C28 and C44. Both these cases are characterized by large-wavelength, wavy features, suggesting an under-representation of sparse roughness in the dataset. Beyond this observation, no clear correlation was found between the error and other primary roughness parameters included herein or surface categorizations (2-D/3-D, random/regular).

The errors associated with using (3.2) are small for surfaces E01 to E06, which were used to calibrate this relation. The errors in using (3.2) and (3.3) over all surfaces in the database are 120% and 430%, respectively. However, when recalibrated against the full database, (3.4) and (3.5) have a significantly smaller error band with maximum values of 79% and 84%. The high error values of the empirical correlations, compared with DNN or GPR prediction, are attributed to the small number of geometrical variables used in their calibrations and the restricted range of the models' parameters.

3.3. Uncertainty estimation

In addition to predictions of equivalent sand-grain height, the GPR method provides confidence margins as functions of each input parameter. These margins can be useful for indicating the kinds of surfaces for which additional training data could improve confidence in predictions. This feature of the GPR approach makes it very attractive for studies of this kind, since DNS and experimental generation of data can be expensive.

The confidence intervals determined by the GPR technique are shown as functions of the normalized surface r.m.s. roughness height, effective slope, porosity and skewness in figure 5. Wider intervals indicate higher estimated values of predictive error, such as at roughness porosity of 0.68, and skewnesses of -1.5 and 2.0 . Surfaces of roughness with similar values of porosity and skewness would then be priorities for additional simulations or experiments.

Equivalent sand-grain height prediction of rough walls

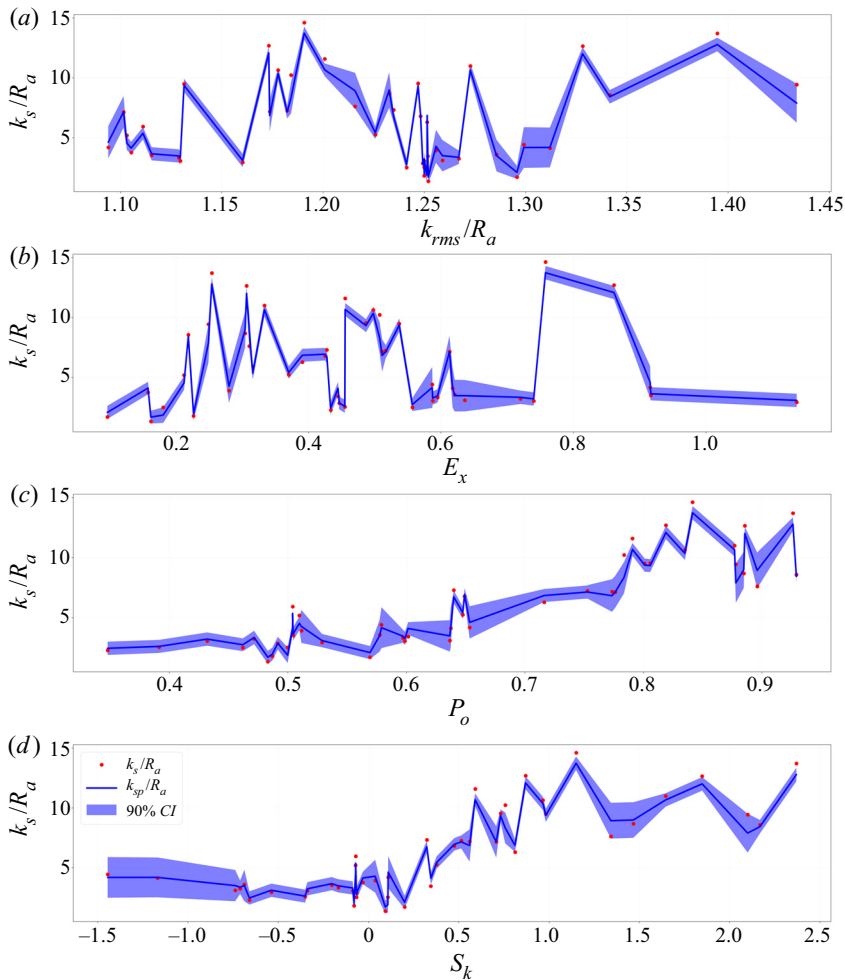


Figure 5. Confidence interval (CI) of predictions with the GPR method, with predicted values of k_s/R_a in blue lines (called k_{sp}) and true values of k_s/R_a in red dots. The GPR predictions for both training and testing data sets are shown – k_s and k_{sp} are very close to each other for the training data points, while they deviate (less than 30 % of error) for some test data points. Line jaggedness is associated with projection of a high-dimensional space to one-dimensional ones.

3.4. Sensitivity analysis

The dependence of DNN predictions of k_s on individual roughness parameters is explored by determining the change in the error norms when each of the primary surface parameters is removed from the data from which the DNN prediction was made. In table 4, the actual error for each surface, and the values of the L_1 and L_∞ norms of errors in the prediction of k_s over the 45 surfaces, are reported when the parameter(s) in the first row is (are) the excluded one(s). The errors of the base prediction (which includes all eight primary parameters) are listed in the second column. In the following discussion, we focus on the L_1 norm for ease of comparison over all 45 cases.

When the values of L_1 are considered, the relative importance of these surface parameters for predicting k_s is E_x , I_x , $|I_z|$, E_z , P_o , k_{rms}/R_a , S_k and, of least importance, K_u . The L_1 -norm error is small when all parameters are included (7.4 %). Excluding any single

Excluded feature(s)	None	E_x	E_z	E_x, E_z	k_{rms}	K_u	k_{rms}, K_u	S_k	P_o	S_k, P_o	I_x	I_z	I_x, I_z
C04	2	-2	3	-1	-1	-2	15	-1	3	13	-12	0	-3
C05	5	-8	11	6	3	-22	-4	0	8	-4	-6	-2	-11
C06	0	10	-1	1	0	10	2	5	1	5	18	6	8
C07	1	3	-1	2	10	-23	0	1	-6	-1	13	1	9
C08	-15	-14	-1	-4	-19	-24	-19	-2	-23	-36	-4	-7	-9
C09	18	4	6	3	0	3	6	1	-2	6	5	11	8
C10	0	1	-16	1	-14	0	-1	-12	2	-13	15	0	11
C11	-12	-3	-3	-23	-2	-2	-5	-12	-1	-29	1	-2	-2
C12	0	-4	-4	0	-18	-3	-4	-1	-7	-2	-3	0	-2
C14	0	4	5	5	1	5	26	3	8	-6	3	6	0
C15	16	5	0	0	2	9	0	0	-11	-2	4	-5	4
C16	1	-2	-1	24	-2	-2	-3	3	-2	6	6	-1	14
C17	-4	8	17	17	1	4	8	15	13	-4	3	5	3
C18	-1	-6	-10	-11	-2	-3	-11	-3	-21	-17	-10	-25	-16
C19	-10	-15	-11	-12	-3	4	5	6	-4	-11	-1	-2	-11
C20	1	3	3	4	3	3	2	4	3	0	23	25	13
C21	9	2	1	3	1	1	2	-1	0	0	0	8	14
C22	-3	-3	-8	-9	-2	-6	-8	-3	-8	-9	-9	-20	-12
C23	0	-2	-1	0	0	-5	-17	-1	0	-1	2	-3	2
C24	0	-21	-1	-1	-1	1	0	0	0	4	0	-4	-7
C26	-6	-17	-12	-9	-8	-5	-19	-15	-13	-5	-13	-14	-10
C28	18	19	21	26	17	18	-3	16	16	32	21	14	20
C29	-9	-19	-8	-22	-6	-5	-13	-25	-11	-22	-18	-17	-19
C30	-4	6	11	25	-10	0	6	24	0	-8	2	6	5
C31	22	20	8	19	24	0	-2	18	-1	-14	9	-1	9
C37	-2	-8	-7	-3	10	-4	-5	-1	-5	-1	-9	-8	-12
C40	-3	-6	-27	-21	-6	-5	-7	0	-1	2	-10	-8	-18
C43	3	-4	-4	6	16	1	2	0	7	23	-15	-1	-12
C44	-6	15	1	17	13	1	4	20	-6	-12	-2	-16	-21
C45	1	2	1	-4	-6	5	-1	-11	1	1	5	2	9
E01	12	4	4	-9	2	-3	-11	5	11	-10	1	-3	-3
E02	-13	6	-6	-7	-2	12	1	-2	10	-9	13	7	-2
E03	15	-6	0	-5	4	-6	-4	3	7	-32	2	1	2
E04	0	-15	-9	-9	-2	-6	-6	-3	-5	2	-2	4	0
E05	5	17	5	17	4	9	9	7	5	28	8	5	13
E06	-5	-3	-6	-3	-10	-9	-10	-6	-7	-9	-10	-10	-5
E07	-21	-21	-24	-18	-16	-21	-18	-17	-23	-41	-25	-25	-24
E08	22	22	25	22	19	18	25	24	7	24	21	22	24
E09	5	-3	15	27	-1	22	26	21	-2	-21	-3	2	2
E10	-18	-19	-5	-8	-25	-4	-5	1	-14	38	-14	8	-2
E11	-1	-15	-23	-19	-7	16	12	-2	9	29	0	-5	0
E12	-9	-3	6	0	-10	2	-2	-15	-10	28	-15	-22	-4
E13	11	8	17	6	17	2	8	7	21	-15	14	25	15
E14	22	6	1	0	6	4	2	1	25	33	9	5	-5
E15	0	18	18	-4	11	9	15	11	19	32	19	23	16
L_1	7.4	8.9	8.2	9.7	7.6	7.1	7.9	7.3	8.0	14.2	8.8	8.6	9.1
L_∞	22	22	27	27	25	24	26	25	25	41	25	25	24

Table 4. Errors in k_s prediction by excluding one or two features. The base prediction includes all primary variables. The four largest errors (in magnitude) for each column are coloured in red. The errors are percentages.

one of these parameters increases the L_1 -norm error up to around 9%. On the other hand, the exclusion of K_u from the input parameters does not worsen predictions of k_s significantly. Instead, this observation appears to be a consequence of correlation between

K_u and other surface parameters like k_{rms}/R_a (see figure 3). When such correlations exist and one correlating parameter is excluded, the DNN process redistributes the weightings given to other correlated parameters, with little loss in predictive accuracy.

To reduce the correlation between the excluded parameters and the remaining ones, one may exclude groups of parameters that are thought to characterize the same type of surface feature. For this reason, a sensitivity analysis was carried out on the effect of groups of variables on predictions of k_s . The characteristics of surface slope, element inclination angle, porosity and intensity of height fluctuations, are contained in pairs of (E_x, E_z) , (I_x, I_z) , (P_o, S_k) and (k_{rms}, K_u) , respectively. Parameters within each pair have been shown to be correlated to some degree in figure 3. Table 4 shows how the accuracy of k_s prediction is affected, if any one of these pairs is excluded. According to the table, the prediction of k_s is sensitive to all four pairs, but with greater sensitivities to the surface porosity (described by P_o, S_k) and the surface slope (described by E_x and E_z). As expected, the elimination of both parameters of a pair worsens the prediction more than removing either single parameter (from around 7–9 % errors to up to 14 %).

According to the sensitivity analysis, all parameters considered are of some importance in the prediction of k_s . The effective x -slope E_x and roughness height skewness S_k have been suggested as especially significant in earlier studies (Napoli *et al.* 2008; Flack & Schultz 2010; Yuan & Piomelli 2014a). The inclination angle in the streamwise direction I_x makes a significant contribution to the k_s prediction because, physically, I_x characterizes the average aerodynamic shape of the roughness elements. Surfaces with $I_x > 0$ are aerodynamically bluff bodies when compared with surfaces of the same size but with $I_x = 0$, and surfaces with $I_x < 0$ tend to be more streamlined and hence produce less drag.

An important finding from this study is that the effective z -slope E_z is of similar importance to accurate k_s prediction as S_k or E_x . The exclusion of E_z adversely affects the prediction for a large number of rough surfaces. Physically, E_z describes whether the surface is close to a 2-D roughness with $E_z = 0$ (such as a transverse bar roughness) or a 3-D roughness with finite E_z . It is known that a k-type 2-D roughness produces a higher drag than a 3-D roughness with the same height due to the larger spanwise length scale that the 2-D roughness imparts to the flow (Volino, Schultz & Flack 2009).

3.5. Comparison between ML algorithms and polynomial models

Explicit algebraic data representations, such as polynomial functions, can also be determined for the data sets of this study, using fitting or minimization procedures. In such methods, a set of basis functions is proposed for a model, the unknown coefficients of which are then optimized according to specified constraints. They are a generalization of the models of equation (1.2a–c), which were based on experimental observations of the dependence of k_s on a small number of surface parameters. A 30-degree-freedom polynomial basis was proposed as a ‘white-box’ model for k_s , analogous to a low-order Taylor series expansion for k_s ,

$$\begin{aligned}
 k_s/R_a = & \alpha_0 + \alpha_1(k_{rms}/R_a)^{\alpha_2} + \alpha_3 I_x + \alpha_4 |I_x|^{\alpha_5} + \alpha_6 |I_z| + \alpha_7 |I_z|^{\alpha_8} \\
 & + \alpha_9 P_o^{\alpha_{10}} + \alpha_{11} E_x^{\alpha_{12}} + \alpha_{13} E_z^{\alpha_{14}} + \alpha_{15} S_k + \alpha_{16} |S_k|^{\alpha_{17}} \\
 & + \alpha_{18} (K_u - 3) + \alpha_{19} |K_u - 3|^{\alpha_{20}} + \alpha_{21} (k_{rms}/R_a)^{\alpha_{22}} P_o^{\alpha_{23}} \\
 & + \alpha_{24} (k_{rms}/R_a)^{\alpha_{25}} E_z^{\alpha_{26}} + \alpha_{27} P_o^{\alpha_{28}} E_z^{\alpha_{29}},
 \end{aligned} \tag{3.6}$$

where a_i ($i = 0, 1, \dots, 29$) are the model coefficients. To keep this model as simple as possible and to bring the effects of all contributing factors into account, we used terms

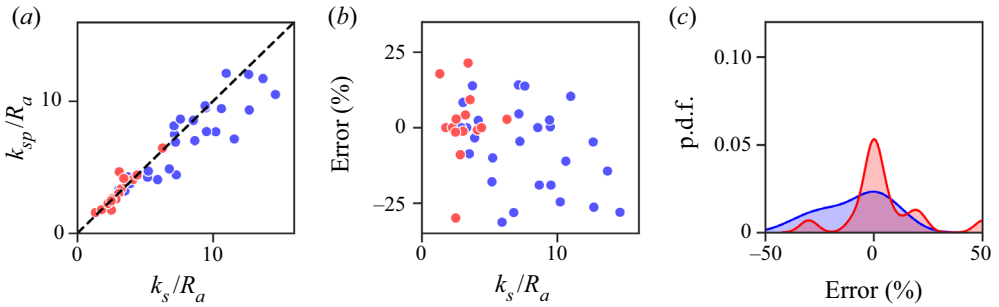


Figure 6. (a) Scatter plot of true k_s and predicted k_s (denoted as k_{sp}), (b) scatter plot of true k_s and relative error and (c) p.d.f. of relative error distribution for prediction using polynomial function defined in (3.6), with DNS data (blue) and experimental data (red).

as $\alpha_i\theta^{\alpha_j}$ for a test variable θ that take only positive values (e.g. k_{rms}), and terms as $\alpha_i\theta + \alpha_j|\theta|^{\alpha_k}$ for those variables that take both positive and negative values (e.g. S_k). For the latter, the power of θ in the first term is fixed (at one) instead of fitted, to eliminate the possibility of an imaginary number. Combinations of six parameters (E_x , E_z , P_o , S_k , k_{rms}/Ra and K_u), taken in pairs, were also included. Since, for the present collection of surfaces, strong correlations were observed between individual variables within the three pairs of (E_x , E_z), (P_o , S_k) and (k_{rms}/Ra , K_u), shown in figure 3, only one variable from each pair was used for the combination terms in (3.6). Using the other variable from any of these pairs instead would not lead to a significant change in the prediction using (3.6).

The high-dimensional space of a_i is poorly suited to curve-fitting and minimization procedures which use stochastic gradient descent algorithms. However, it is well suited to robust minimization methods like the differential evolution algorithm (Storn & Price 1997), with which global minima can often be found efficiently in spaces of high dimension. In this case, it is used to determine the values of the coefficients a_i which minimize the L_1 norm.

In figure 6, the prediction quality of this white-box model with optimized coefficient values is shown. This method yields an average prediction error of 12 % and a maximum one of 51 % when using all 45 fully rough data sets (to give the best possible prediction accuracy) for the model training.

The optimized values of a_i are

$\alpha_0 = 5.312,$	$\alpha_1 = -1.172,$	$\alpha_2 = 4.264,$	$\alpha_3 = 0.050,$	$\alpha_4 = -1.283,$	$\alpha_5 = 8.393,$
$\alpha_6 = -0.347,$	$\alpha_7 = -5.771,$	$\alpha_8 = 1.785,$	$\alpha_9 = 7.919,$	$\alpha_{10} = 4.058,$	$\alpha_{11} = -0.979,$
$\alpha_{12} = 3.414,$	$\alpha_{13} = 6.380,$	$\alpha_{14} = 1.354,$	$\alpha_{15} = 1.023,$	$\alpha_{16} = 2.969,$	$\alpha_{17} = 1.273,$
$\alpha_{18} = -0.946,$	$\alpha_{19} = -0.762,$	$\alpha_{20} = 0.056,$	$\alpha_{21} = 1.647,$	$\alpha_{22} = -8.176,$	$\alpha_{23} = 3.523,$
$\alpha_{24} = -9.472,$	$\alpha_{25} = -5.656,$	$\alpha_{26} = 0.580,$	$\alpha_{27} = -5.425,$	$\alpha_{28} = 0.283,$	$\alpha_{29} = 7.177.$

The predictive accuracy of this optimized explicit model equation is considerably lower than that of the DNN and GPR methods. One reason for this reduced accuracy is that low-order functions of geometrical parameters do not faithfully represent the dependence of k_s on surface parameters because each coefficient in the model is required to take the same value over the entire surface-parameter space. In ML approaches, such restrictions need not apply as they are not constrained to low-order polynomial functions but instead adopt a methodical search for the best representation of k_s as a function of the surface

parameters. This search is carried out through ‘feature selection’ in the first layers of DNN and the properties of the basis functions adopted in GPR, each of which are designed to yield the same mean and standard deviation of k_s/R_a as in the original dataset (Rasmussen & Williams 2006).

4. Concluding remarks

The construction of a predictive model from a large ensemble of datasets for the equivalent sand-grain height k_s of a surface of arbitrary roughness, as a function of many different measures of surface topography, is a labelled regression problem that is well-suited to ML techniques. In this paper, data from 45 different rough surfaces (in fully rough flows) were used to devise DNN and GPR predictions for k_s as functions of eight different surface-roughness parameters.

Both models were able to predict k_s for the 45 surfaces with an average error below 10 %, with the largest error for any one surface less than 30 %. These predictions were significantly better than those of existing formulae, and of a 30 degree-of-freedom polynomial model fitted to the same data, where the greatest error for any surface was approximately 50 %.

Sensitivity analyses revealed that inclusion of nearly all the surface roughness descriptive parameters was necessary to minimize the average prediction error, but that exclusion of either measures of porosity or measures of the surface slope increased the maximum prediction error more significantly than omitting other parameters.

Machine learning techniques are well suited to this modelling problem because: (i) it is complex in so far as different kinds of surface roughness yield different flow phenomena which are modelled most accurately in different ways, making the prospect of a general physical model very remote; and (ii) the dependent surface-roughness variables upon which k_s is modelled are a large non-orthogonal set for which robust multivariable regression techniques are required. As ML methods, they take no account of physical modelling concepts or observed phenomena within roughness sublayers, such as recirculation regions, enhanced turbulence production in the wake of roughness elements, assumed scalings for drag, etc., each of which is applicable to flows over some rough surfaces but not others. Nor are they hindered by the lack of orthogonality of the surface roughness parameters as the dependent variables of k_s . The techniques used can be configured readily to mimic models with very many degrees of freedom and, when compared with polynomial models, their feature selection properties provide the equivalent of different values for polynomial coefficients in different regions of the surface-parameter space. In this application, both approaches of DNN and GPR yielded models with very similar predictive accuracy, even though the techniques themselves were very different. We therefore conclude that they yield high-fidelity predictions of the equivalent sand-grain roughness height for turbulent flows over a wide range of rough surfaces, as a significant improvement over other methods. Improved prediction might be achieved by enlarging the database to include rough-wall flows with surface parameters which correspond to the relatively low prediction confidence in the GPR method, and by including additional roughness parameters as inputs which might describe sparseness and two-dimensionality, such as the solidity, correlation length scales and other two-point surface statistics.

In addition to the k_s prediction described here, the DNS database and the ML techniques in general can also be used to uncover relations between roughness geometry and physics-related quantities, such as the flow pattern around roughness protuberances, flow

separation locations, characteristics of the shear layers associated with the separation bubbles, the wake sheltering volume, etc. Specifically, an ML network trained to correlate these flow characteristics (as outputs) to the roughness geometry (as inputs) may be an efficient tool for determining the sets of roughness geometrical features which are important for characterizing these effects. Knowledge of such a set of significant roughness parameters may also guide the construction of rough-surface databases that yield more efficient and more widely applicable predictions of k_s or other quantities.

Supplementary materials. The rough-wall flow database (including k_s , surface height map and surface parameters) and the trained DNN and GPR networks, called prediction of the roughness equivalent sand-grain height (PRESH), can be accessed online in the first author's GitHub repository at <https://github.com/MostafaAghaei/Prediction-of-the-roughness-equivalent-sandgrain-height>. With this package of data and programs, interested researchers can: (i) use the ML networks described in this paper to make predictions of k_s for surfaces of their own roughness topography; (ii) download the code and train new DNN and GPR networks to predict k_s for a different set of surfaces of arbitrary topography; and (iii) use the database of 45 rough-wall flows for other applications. It is recommended to use the ML configurations described in this paper for surfaces with parameters inside the ranges specified in figure 3. Extrapolations (using inputs which are beyond the specified range) will lead to additional uncertainty.





The PRESH and the database will be actively updated by the authors to improve the prediction accuracy and universality. We welcome interested researchers to share their datasets with us.

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