The authors would like to thank the reviewer for the comment. The novelty of the present study is two-fold, as mentioned below:

1. The force partitioning method allows the decomposition of total force into kinematic, vortex-induced and viscous forces, which helps to understand the driving force during the initiation and sustenance of VIV. The implementation of the methodology in OpenFOAM is the first of its kind, which is explained in the manuscript.
2. Contrary to the work of Viré et al. [4], the force partitioning method is used to analyse the force contribution during VIV at transcritical Reynolds numbers. Tower properties of IEA 15MW are used for the study, which was not the case in the previous literature.

According to the previous works of Belloli et al., [1], the surface roughness may be used to mimic the high Reynolds number phenomena at a considerably lower Reynolds number. This methodology can shed light on the contribution of surface roughness on VIV at lower Reynolds numbers in the future.

Most of the above-mentioned comment is explained in the last paragraph of the introduction section. However, to emphasize the novelty of this work further, additional text is added to the conclusion.

Line 80: While the previous study focuses on the flow around a cylinder at Reynolds number $\leq 250$, the importance of force partitioning method at transcritical Reynolds numbers is yet to be understood, as the flow around wind turbines is characterised by a high Reynolds number ($Re > 3.5 \times 10^6$).

Line 390: Even though force-partitioning analyses have been carried out in previous studies, investigating the phenomenon of VIV at such high Reynolds numbers with this method is the first of its kind.

2. The rationale behind the 3D-$\rightarrow$2D casting should be further illustrated in the manuscript. The authors have selected a section corresponding to the tower top. Is it the most critical section with regards to a potential lock-in?

As a wind turbine tower is a tapered structure, analysis of such a three-dimensional structure in two-dimensional is not straightforward. The purpose of simplifying the analysis is to study the force partitioning methodology on a simple geometry at high Reynolds numbers. Future studies will focus on numerical studies of complete three-dimensional towers. It is also to be noted that the simulations carried out in OpenFOAM are never completely two-dimensional, but the mesh is extruded with a single cell thickness in the third dimension.

The structural properties, such as mass and stiffness, vary along the height of a three-dimensional tower. Moreover, a wind turbine tower can be considered as a cantilever beam standing perpendicular to the ground. Hence, the bending characteristics of such a structure during VIV cannot be obtained in a two-dimensional study. To simplify the overall analysis, the section corresponding to the tower top is used for the present study, as the overall displacement is maximum at the tower top and not due to being a critical section in regards to a lock-in. The following changes are made to lines 265-268.

Line 265-268: The characteristics of the top section of the IEA 15MW reference wind turbine are chosen as the analysis of a three-dimensional structure in two-dimensional is not straightforward, and it simplifies the overall analysis. The tower top section is chosen as it experiences the maximum displacement in the scenario of VIV and is not due to being a critical section during a lock-in.

3. This reviewer found problems seeing how the values of Table 2 ($m^*, c^*$) were computed,
and what their relation is to the mechanical properties and dimensions of the full tower.

The mass ratio is calculated for the whole tower as shown below.

\[ m^* = \frac{m_L}{(1/4) \times \pi \times \rho \times D^2}, \quad (1) \]

where \( m_L \) is the mass per unit length, \( \rho \) is the density of the air and \( D \) is the average diameter considering the base and top. The mass ratio is calculated for the top section of the tower with a height of 12.386 m. The authors would like to thank the reviewer for identifying the typographical mistake in the value, which has been rectified to 87.6. With the consideration from another reviewer, \( c^* \) is changed to \( \zeta \) as explained in line 167. The damping for the tower is 2\%, and the damping ratio is calculated with the following equation.

\[ \zeta = \frac{\delta}{\sqrt{4\pi^2 + \delta^2}}, \quad (2) \]

where \( \delta \) is logarithmic decrement.

4. Regarding the FSI implementation, did the authors employ a weak or a strong coupling approach? How was the mesh deformed?

In the present study, a strong coupling algorithm is used for both flows in the laminar and turbulent regime. The mesh deformation is carried out with Spherical Linear Interpolation Scheme (SLERP) interpolation which is already implemented in OpenFOAM.

Line 201: A strong coupling algorithm is utilised for the present study to give accurate results for modeling VIV. The structural parameters are computed for each iteration of the fluid solver, where the fluid solvers are sub-iterated for multiple iterations till the convergence criteria are met. The computed displacement of the cylinder moves the mesh according to Spherical Linear Interpolation Scheme (SLERP).

5. Consider including a validation for the turbulent setup, or citing Viré et al. (2020) to that end.

As explained in Section 3.2 of the manuscript, the grid-independent study is compared against the results observed by Viré et al. [4], where a similar simulation setup and study were performed. The present work now cites the previous work as extensive validation and a comparative study is made in the previous work. The following sentence is added to line 289, explaining the same.

Line 289: As explained in Section 3.2, the simulations performed at Reynolds number, \( Re = 3.6 \times 10^6 \), serve as a validation for the simulations at the turbulent regime. The force coefficient, pressure distribution, and separation point align with the results observed from the work of Viré et al. [4].

6. Equation 10: The amplitude factor \( y_0 \) was never introduced. Even if not relevant for this work or for the derivation, the generalized form of the equation may also take an initial phase. I found the comment ”The mode shape at any time represents the ratio of the displacement at the point \( z \) to the reference displacement \( y(t) \),” rather confusing. The mode
shape is time-independent, and the reference displacement has not been introduced. The choice of axes notation is also confusing, as e.g. the axial direction of Figure 1 corresponds to “x”.

The authors understand the confusion derived from the paragraph and the changes for lines 177-179 are made as shown below. Figure 1 is edited for better understandability.

Line 177: where $\Psi(z)$ is the shape function, which represents the ratio of the displacement at the point $z$ to the reference displacement $y(t)$. The reference displacement $y(t)$ is represented as a sinusoidal function with amplitude $y_0$.

7. Statistics shown in Figure 8 may suffer from the short total times simulated. Could be interesting to include a comment in the manuscript, and to state if that could be a limitation for the analysis made.

The authors would like to thank the reviewer for the comment, and the following text in red is added to line 315 of the manuscript. As an explanation, the simulations were stopped when the vorticity-induced force became fully transient in nature. The simulations conducted were free vibration, and the oscillations can reach even higher if there is insufficient negative aerodynamic damping. Due to a considerable amount of time, cost and the transient nature of the vorticity-induced force, the simulations were stopped at this particular behaviour.

Line 315: As seen in Figure 9, the simulations were stopped once the vorticity-induced force became completely transient in nature. The system becomes self-exciting and the fluid imparts energy to the structure if the structural damping coefficient is not enough to suppress the aerodynamic excitement. This can be analysed with a longer simulation time, which is out of scope for the current study due to increased time and cost.

8. In Figure 12, for the moving average, was the considered window a multiple of the period of motion? In the opinion of this reviewer, that could facilitate future comparisons.

The non-dimensional energy ($E^*$) extracted by the cylinder is calculated according to equation 24 of the manuscript. The mean of $E^*$ is then calculated for every period of the energy cycle, which is a multiple of the period of force fluctuation.

9. ”As the oscillations are almost sinusoidal, the net energy per cycle becomes almost zero.” This reviewer had problems understanding this statement. Even with pure sinusoids, a phase difference should lead to a non-zero energy transfer (when considering what happens at the frequency of motion). Note that, at the LCO of e.g. Figure 9.a, one should be ”pumping” energy into the system to counteract the structural damping and maintain the same amplitude of vibration.

The authors understand the confusion created by the sentence, and it is struck out to avoid further doubts. During each energy cycle, the mean non-dimensionalised energy transfer from kinematic force is negligible when compared to that of vorticity-induced force. Especially during the initiation of oscillations, much of the energy transfer happens from fluid to structure due to the vortices formed behind the cylinder. The
total force experienced by the cylinder is largely driven by this, which can also be seen in Figure 9c,e. Once the structural damping overcomes the forces derived due to vortices, the total force becomes out of phase with the displacement and beating phenomenon, as in Figure 9c, can be seen.

As the oscillations are almost sinusoidal, the net energy per cycle becomes almost zero. Hence any transient behaviour or sudden change in lift force comes purely from the vorticity-induced force.

10. While the distribution of the OpenFOAM implementation performed by the authors is highly appreciated, and it is well described in the manuscript, it could be interesting to improve the writing so that non-OpenFOAM practitioners do not get lost. For instance, one could include more ”fundamental” descriptions of what the ”0” folder contains, or how blockMesh generates the grids.

In order to keep the manuscript related to the methodology and its implementation, the fundamental explanations of running OpenFOAM simulations are now updated in the readme of the GitHub repository. This is now explained in the manuscript in line 142.

Line 142: The source code of the implementation can be found in the repository, where the fundamental description of running the OpenFOAM simulation is explained in detail.

Minor Comments

11. Could be interesting to comment on how the method should be extended to situations where the structure is characterized by more than one mode.

It is expected to carry out such studies in the future with multiple structural modes and environmental conditions. As the simulations in OpenFOAM are not completely two-dimensional, the methodology or the implementation does not change. However, carrying out high-fidelity or mid-fidelity simulations can give an insight into which part of the structure gets excited and initiates the VIV, and vice-versa, and how VIV are damped. Breaking down the domain into different parts of the fluid domain also gives an insight into how the near wake and far wake regions affect the VIV, as can be seen from the studies of Menon et al. [3].

12. Is ”Cartesian grid” the best way to describe the mesh? I believe near the cylinder your grid lines become normal to the surface.

It is now renamed to O-grid topology in line 147.

Line 147: As it is a relatively simple geometry, a body-fitted Cartesian grid mesh with O-grid topology is made to carry out the simulations.

13. ”Nevertheless, the more recent study by Viré et al. (2020) gives confidence in performing URANS simulations”: while this is true, could passing to another turbulence model eventually have an impact on the conclusions of this work, for instance regarding the relative contribution of the forces?
The present study uses \( k-\omega \) SST as the turbulence model, proving to predict fluid separation and strong pressure gradients better. The present work matches well with the results from the previous experimental and numerical studies, which is the primary reason for choosing it as the turbulence model. Another turbulence model like \( k-\epsilon \) seems to underpredict the skin friction with a similar boundary condition and is known to perform poorly on flows with large separation. The flow is found to have a delayed separation and lower base pressure, which eventually gave slightly higher drag as per the studies by Catalano et al. [2].

The physical parameters analysed from the numerical studies with different turbulence models can have different results for pressure coefficient, separation point, wakes, etc., which are inherent to the turbulence modelling itself. The relative contribution of forces could then vary depending on the above-mentioned physical parameters. However, the overall calculation of force, depending on the direction in which it is analysed, will be the same as global force parameters like lift and drag.

14. Would the provided OpenFOAM modifications work for 3D cases in its present form?

OpenFOAM is a code defined for three-dimensional space and the meshes are defined as such. To carry out a two-dimensional analysis, the mesh is made with a single-cell thickness, and suitable boundary conditions are provided at the patches that are normal to the direction of interest. Therefore, the code works perfectly for 3D cases in the present form.

15. The link to Zenodo of the manuscript did not seem to be correct (but a working one was indicated by the authors in the submission).

The URL is now changed to a working link.


16. Equation 9: the symbol used here for the natural frequency is later on used in the manuscript for displacements.

The variable is changed from \( \nu \) to \( v \) for the equation 10, 11 and 12.

17. Equation 11: for consistency with the rest of the derivation, maybe it is a good idea to use ”’ for the second derivative.

The equation is changed to maintain consistency as recommended.

18. Equation 16: there is a jump in the derivation here, from continuous to discrete. Was it intended?

As equation 14 explains, the natural frequency is obtained by equating potential energy and kinetic energy. On substituting the shape function to this equation, equation 16 can be obtained. One thing to consider here is that the differentiation term \( \Psi'' \) is already carried out and the final form is written in equation 16.

19. ”frequency of the tower is calculated to be 0.48” -> 0.48 Hz
The correction is now implemented.

20. **Figure 3: misses the coordinates for the center of the cylinder.**

Figure 3 is changed to include the upstream and downstream size of the domain, which is sufficient to determine the location of the cylinder in the domain.

21. **Some words on the computational cost of the simulations would be appreciated.**

The simulations are carried out with 20 cores, where one of the most expensive simulations required approximately 4950 CPU hours, for a velocity ratio of $U/U_{st} = 0.997$, and 2520 CPU hours, for $U/U_{st} = 0.71$, using Intel(R) Xeon(R) CPU E5-2670v2 processor. This is added to line 318.

References


Reviewer 2

Specific Comments

1. The ”subcritical” Reynolds number regime is generally understood to be different than the laminar Reynolds number regime (see for example the cited work by Belloli (2015)). In the present work, both terms are used interchangeably.

In the present work, the simulations performed at subcritical Reynolds numbers are between 80 and 200. As the Reynolds numbers are low enough to form a laminar flow around the cylinder, the authors have mentioned it as a laminar Reynolds number regime. But, for a generalised term where the flow is no longer laminar (especially for Re > 250) but subcritical Reynolds number range, the term subcritical is used in the manuscript.

2. Chapter 3.1: The reviewer is missing boundary conditions at the inlet for $k$ and $\omega$. There is information about $k$ and $\omega$ in lines 256ff but they are, at least for $k$, introduced as initial values.

The authors would like to thank the reviewer for the comment. The following sentences are added to the manuscript at line 286.

Line 286: The boundary conditions for $k$ and $\omega$ are Dirichlet boundary conditions at the inlet with values equal to initial values as calculated above. Neumann boundary condition is imposed for both the variables at the outlet.

3. Chapter 3.1: The reviewer is interested in the time steps used and the resultant Courant number. Has the used time step been verified?

An adaptable time step is adopted for the simulations performed in the present study. A maximum Courant number of 0.7 is set to determine the time step for each time iteration, which is verified in the previous study by Derksen [3].

Line 234: The time step used for all the simulations maintains Courant–Friedrichs–Lewy (CFL) number equal to 0.7.

4. Chapter 3.1: The reviewer is interested in the coupling approach of fluid and body.

In the present study, a strong coupling algorithm is used for both flows in the laminar and turbulent regime. The following sentence is added to line 201, explaining the coupling algorithm.

Line 201: A strong coupling algorithm is utilised for the present study to give accurate results for modeling VIV. The structural parameters are computed for each iteration of the fluid solver, where the fluid solvers are sub-iterated for multiple iterations till the convergence criteria are met. The computed displacement of the cylinder moves the mesh according to Spherical Linear Interpolation Scheme (SLERP).

5. Figure 3/numerical domain: Information about the size of the upstream and downstream
part of the domain is missing.

Figure 3 is changed to include the upstream and downstream size of the domain.

6. Figure 3/numerical domain: The domain in the across-wind direction seems rather small with 15 d and a blockage ratio that is above 7%. The cited work by Viré (2020) uses 100 d in the across-wind and 100 d in the along-wind direction. Verification of the domain size can show the appropriateness of the chosen domain size.

The blockage ratio from the current domain size is 6.67% (Frontal area of cylinder/Inlet area of the domain = \( \frac{d}{15d} \)). Even though the work by Viré et al. [6] has adopted 100d along-wind and across-wind direction, the domain size can be reduced for quicker calculations with more mesh points, without having reflections from the domain boundary. As mentioned in the literature study by Rodriguez et al. [5], the current blockage ratio falls within the values mentioned in the previous studies on VIV. Moreover, the separation point or transition point will not be affected as the wake width decreases as Reynolds numbers reach supercritical Reynolds number ranges [1].

The simulation is validated and verified for subcritical and transcritical Reynolds number regimes as explained in Section 4.1 and 3.2 respectively. The numerical simulations at Reynolds number 3.6 \times 10^6 is validated against the work of Viré et al. [6], where satisfactory accuracy is obtained as mentioned in Section 3.2.

7. Chapter 3.2: The grid convergence study is conducted for the surface pressures on a fixed cylinder in transcritical Reynolds number regime. The main aim of the paper, however, is to investigate the situations for a SDOF oscillator. In the reviewer’s view, the grid convergence study should investigate a critical situation during free vibration. Additionally, the increase of number of cells by a factor of 2 seems rather small even for a 2D simulation.

The authors believe that the Grid Convergence Index (GCI) conducted for surface pressures on a fixed cylinder is still scientifically relevant in the present study. While investigating the situations for a single degree of freedom oscillator, another complexity of mesh motion adds up in the GCI study, which cannot be separated from the discretisation error. Nevertheless, the comparison of mean surface pressure on a fixed cylinder not only confirms the global integrated variables like force coefficients but also the surface pressure at every point around the cylinder. Moreover, the validation done at subcritical and transcritical Reynolds numbers adds to the confidence of the result.

According to the procedure estimated by Celik et al. [2] and Roache [4], Richardson’s extrapolation technique is used to estimate GCI. A grid refinement factor greater than 1.3 is suggested by the literature, which can be attained by doubling the number of mesh points. In the present study, the grid refinement factor is approximately 1.413.

8. Table 1: The reviewer is interested which \( y^+ \) value is shown. Is it the mean value?

This is the value at which the first cell height of the mesh is calculated. The first cell height is kept constant for all the meshes, while the mesh growth ratio is changed.
9. line 209f: How is the separation point calculated and where is it located?

The point where the wall shear stress becomes zero is taken as the separation point on the cylinder. The separation points for three meshes are as follows: Coarse - 112.9°, Medium - 112.2°, and Fine - 111.8°. The line 222 is modified as shown below.

Line 222: Furthermore, the medium and fine mesh predict the separation point to be 112.2° and 111.8° respectively, in comparison to 111° from the literature.

10. Figure 5, 6, and 8: It is not clear how the statistical values are obtained. Is the whole time history used, or is an initial time of the simulation discarded?

The mean drag coefficient and r.m.s of lift coefficient are calculated after the simulation achieved a steady state. However, the whole time series is considered for the calculation of the maximum lift coefficient and maximum non-dimensionalised amplitude of oscillation.

11. Figure 5, 6, and 8: The reviewer is missing information about the verification of the convergence of the results. It is not clear how the chosen simulation time has been deemed appropriate to calculate the statistical values.

The results obtained for flow in the laminar regime are concluded after developing steady-state oscillations. Hence the verification of convergence of results is considered reliable for all the simulations in this Reynolds number regime. But for simulations carried out at turbulent regimes, the simulations were stopped after vortex-induced force became fully transient. As the simulations performed were free vibration, the structural damping was insufficient to restrict the oscillation at this point in time, especially around the lock-in region. In all the other velocity ratios, especially when the oscillations were not as significant as lock-in, harmonic steady-state solutions were obtained. The following sentence is now added, line 315, explaining the same for flows in turbulent regimes.

Line 315: As seen in Figure 9, the simulations were stopped once the vorticity-induced force became completely transient in nature. The system becomes self-exciting and the fluid imparts energy to the structure if the structural damping coefficient is not enough to suppress the aerodynamic excitement. This can be analysed with a longer simulation time, which is out of scope for the current study due to increased time and cost.

12. One of the main concerns of the reviewer are the already mentioned diverging forces in Figures 9 c) and e). To the best of the reviewer’s knowledge, the lift force does not diverge during VIV. See for instance the wind tunnel results by the co-author Belloli (2012) and Belloli (2015). In the numerical simulations of the co-author Viré (2020) it is shown that the lift force decreases significantly after reaching a maximum in Figure 11 $\omega^* = 0.95$. Is this change of lift coefficient in time a particular situation that only appears in URANS simulations or has it been already observed in wind tunnel or full-scale?

As seen from Zasso et al. [7], there is an irregular development of lift coefficient at the beginning of the lock-in region, where a phase difference much greater than zero can be seen. After a certain time, the phase difference drops to almost zero, and the lift coefficient and the oscillation amplitude are perfectly
synchronized. However, this phenomenon is also seen in URANS simulations because there is a switch in phase difference from positive to negative. This is mainly due to the change in wake pattern, which eventually dampens the oscillation, which causes the force to shift from a self-exciting to a self-limiting effect [3]. This can also be seen in the present study with the phase difference between kinematic and vorticity-induced force, as explained in section 4.2.1 of the manuscript.

13. The reviewer is interested to know if the oscillation amplitudes also diverge. VIV are self-limited and the oscillation amplitudes should reach a maximum even during lock-in, contrary to aeroelastic phenomena such as galloping or flutter.

As seen from Figure 8, VIV is a self-limiting phenomenon where large oscillations occur near the lock-in region. Any higher or lower wind speeds result in little to no vibrations in the structure. In the current study, the maximum amplitude is reached at a velocity ratio of 1 at a Reynolds number of $4.6 \times 10^6$.

14. Figures 9 c) and e): The reviewer is missing information about the criterion used to stop the simulations after 450 s and 525 s, respectively. What happens after the shown time histories?

As explained in comment 11, the simulations were restricted as the vortex-induced force reached fully transient. The structural damping was insufficient to limit the oscillation at this point of time for simulations in the lock-in region. The simulations are also stopped as the computational time and cost became considerably high and the uncertainty of not knowing when an equilibrium is reached. The following sentence, as explained in comment 11, is added in line 315.

Line 315: As seen in Figure 9, the simulations were stopped once the vorticity-induced force became completely transient in nature. The system becomes self-exciting and the fluid imparts energy to the structure if the structural damping coefficient is not enough to suppress the aerodynamic excitement. This can be analysed with a longer simulation time, which is out of scope for the current study due to increased time and cost.

Technical corrections

15. line 22f: ”challenge” instead of challenges, ”as taller” instead of ”as a taller”

The correction is made in the manuscript.

16. Figure 4: The definition of the normalized x-coordinate $x^*$ is missing.

The definition for normalised x-coordinate, as shown below, is explained in line 217.

Line 217: The normalised x-coordinate $x^*$ is the x-coordinate normalised to the chord of the cylinder, as $x/c$.

17. The direction of lift and drag and the respective formulas for the coefficients could be defined for clarification.
The lift and drag forces are defined in line 239 at the beginning of the results. The equations for lift and drag are shown in Equation 19.

Line 239: The lift and drag coefficients calculated from the OpenFOAM postprocessing function tool are explained in Equation 19, where $u$ and $v$ are the velocity components in the x-direction and y-direction, respectively. The lift and drag force are defined as $F_L$ and $F_D$, respectively, and the surface area of the cylinder as $S$.

18. Figure 5 caption: remove "citep".

The caption is edited for correction.

19. line 236 "y-direction correponding" add space after comma.

The correction is now implemented.

20. Table 2: Damping ratio has been introduced as $\zeta$ on p.8 instead of $\zeta^*$ here. Is there a difference between these two damping ratios?

It was a typographical error from the authors and it is now changed to $\zeta$.

21. line 264: $y_{wall}$ is not defined.

The first cell height of the mesh is defined by the variable $y_{wall}$, which is defined in line 2.


The URL is now changed to a working link.


References


Reviewer 3

Remarks

1. Within the text, the force related to shedding is called, alternatively, vorticity force and vortex-induced force. The authors should uniformise how they refer to it.

The authors confirm the non-uniformity, and the word has been replaced with vortex-induced force throughout the manuscript.

2. The added mass force, first term in the RHS of equation 4, is never properly defined. Moreover, while maybe obvious, the difference between the volume V and the surface B in integrals is not given in the text.

The first term in equation 4 corresponds to the added mass effect. It is determined by the rate of change of velocity of the object. Hence, it is purely dependent on the velocity and shape of the object, as it is integrated over the surface B. The equation is briefly explained in lines 110-121, where the detailed explanation can be found in Menon et al. [2], as mentioned in line 111. The terms V as volume and B as surface are explained in lines 98 and 99. The equation is now changed to maintain consistency in the terms for volume from V to \( V_f \) as shown below.

\[
F_B = F_K + F_\omega + F_\phi + F_\varepsilon
\]

\[
F_B = -\rho \int_B \vec{n} \cdot \frac{d\vec{U}_B}{dt} \phi^{(i)} dS - \rho \int_B \frac{1}{2} \vec{U}_B^2 \vec{n} \cdot \nabla \phi^{(i)} dS + \rho \int_{V_f} \nabla \cdot (\vec{\omega} \times \vec{n}) \phi^{(i)} dV + \rho \int_{V_f} \nabla^2 \left( \frac{1}{2} \vec{u}_\nu^2 + \vec{u}_\nu \cdot \vec{u}_\nu \right) \phi^{(i)} dV
\]

\[
+ \mu \int_B (\vec{\omega} \times \vec{n}) \cdot \vec{\nabla} \phi dS + \rho \int_{V_f} \nabla^2 \left( \frac{1}{2} \vec{u}_\nu^2 \right) \phi^{(i)} dV - \rho \int_{\Sigma} \frac{d\vec{U}}{dt} \cdot \vec{n} \phi^{(i)} dS + \mu \int_{\Sigma} (\vec{\omega} \times \vec{n}) \cdot \nabla \phi^{(i)} dS
\]

3. Figures 1 and 3 are inconsistent. Moreover, if the cylinder is 2D, what is the height H in table 1? It would help to understand this parameter if it is added to figure 3.

Figure 1 is changed to make it consistent with figure 3. It is to be noted that Figure 1 is more representative in nature and not the domain characteristics of the present study. The domain characteristics explained in Figure 3 are shown as two-dimensional, as the mesh extrudes with a single cell thickness, unlike a fullfledged three-dimensional mesh or numerical study. It is carried out in such a way as OpenFOAM is a code designed for three-dimensional space and cannot have two-dimensional mesh. The boundary conditions in the patches normal to the direction of interest are made to make the flow completely two-dimensional. This is explained in lines 162 and 273.

Line 162: As OpenFOAM is a code designed for three-dimensional space, the mesh is extruded in the z-direction with one cell thickness.
Line 273: The parameter \textit{height} shows the single cell thickness in the z-direction as mentioned in Section 3.1.

4. It is not clear which criteria were used to choose the set of parameters (c, EI, ks) from equations 9-18 (or the mass ratio in table 2).

The structural parameters mentioned in equations 9-18 are calculated as follows:

Youngs modulus, E for IEA 15MW tower = \(2 \times 10^{11}\) N/m\(^2\)

Moment of inertia for each section is calculated as

\[ I_j = \frac{\pi (D_{out}^4 - (D_{out} - D_{in})^4)}{64}, \]  

where \(I_j\) is the moment of inertia of a section, \(D_{out}\) is the outer diameter of the cylinder, \(D_{in}\) is the inner diameter of the cylinder. The equivalent moment of inertia is calculated as mentioned in equation 17 in the manuscript.

Spring constant, \(k\) is calculated from the equation

\[ k = m\omega^2, \]  

where \(m\) is the mass of the tower/cylinder.

The parameters of the tower were chosen from the IEA 15 MW report as per IEA Task 37 [1]. The mass of the cylinder with a height of 1 m is calculated from the mass ratio of the IEA 15 MW tower. Equation 2 mentioned above is added to line 193 in the manuscript and the equation 3 is mentioned in line 167.

Line 193: where the moment of inertia of a section of the tower is calculated as

\[ I_j = \frac{\pi (D_{out}^4 - (D_{out} - D_{in})^4)}{64}, \]  

and \(D_{out}\) is the outer diameter of the cylinder, and \(D_{in}\) is the inner diameter of the cylinder.

5. Why no comparison with the literature has been made for the cylinder in the turbulent regime? This could be done, at least, for a non-oscillating system. This would be still relevant as the turbulent wake may be conditioned by the turbulence model, separation point, etc.

A grid-independent study is carried out in the present study, where the force coefficient, pressure distribution and its difference are simulated for various meshes. As seen in Section 3.2 of the manuscript, the results are compared with the works of Viré et al. [3], where a similar simulation setup and study was performed. An extensive validation and comparative study was made by Viré et al., and the present work is referenced to their works, as the results align with the previous study. The following sentence is added to line 289, explaining the same.

Line 289: As explained in Section 3.2, the simulations performed at Reynolds number, \(Re = 3.6 \times 10^6\), serve as a validation for the simulations at the turbulent regime. The force coefficient, pressure distribution, and
separation point align with the results observed from the work of Viré at al. [3].

6. Overall, the labels in figures are too small and hard to read even in electronic form.

The figures are made bigger for better legibility.

References

