Nektar++: Development of the Compressible Flow Solver for Large Scale Aeroacoustic Applications

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Abstract

A recently developed computational framework for jet noise predictions is presented. The framework consists of two main components, focusing on source prediction and noise propagation. To compute the noise sources, the turbulent jet is simulated using the compressible flow solver implemented in the open-source spectral/hp element framework Nektar++, which solves the unfiltered Navier-Stokes equations on unstructured grids using the highorder discontinuous Galerkin method. This allows high-order accuracy to be achieved on unstructured grids, which in turn is important in order to accurately simulate industrially relevant geometries. For noise propagation, the Ffowcs Williams - Hawkings method is used to propagate the noise between the jet and the far-field. The paper provides a detailed description of the computational framework, including how the different components fit together and how to use them. To demonstrate the framework, two configurations of a single stream subsonic jet are considered. In the first configuration, the jet is

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treated in isolation, whereas in the second configuration, it is installed under a wing. The aerodynamic results for these two jets show strong agreement with experimental data, while some discrepancies are observed in the acoustic results, which are discussed. In addition to this, we demonstrate close to linear scaling beyond 100,000 processors on the ARCHER2 supercomputer. *Keywords:*

Jet Noise; Discontinuous Galerkin; Scaling; Large Eddy Simulation; Ffowcs Williams - Hawkings; Acoustics; Installation Noise.

1 1. Introduction

Many fluid flows of engineering interest generate noise, either directly, 2 or indirectly through interactions with nearby solid surfaces. Examples of 3 this include the flow around cooling fans in computers [1], the flow around A-4 pillars and rear view mirrors on cars and trucks [2], the flow over wind turbine 5 blades [3], the flow around the undercarriages of high-speed trains [4], and the 6 flows generated by different parts of an aircraft, including high-lift systems [5, 6], fan- and turbine blades [7, 8], landing gears [5, 6], and the propulsive 8 jet [9, 10, 11]. This type of aerodynamically generated noise will often be 9 perceived as unpleasant, and can in some cases even be harmful. Therefore, 10 there is a widespread need to understand and minimize the underlying noise 11 sources. 12

The study of aerodynamically generated noise, known as aeroacoustics, saw a great surge in the middle of the last century. In the beginning, the motivation for studying aerodynamically generated noise came mainly from the aircraft industry, where it was realized that jetliners equipped with the

newly invented turbojet engine would be too loud for widespread use. This 17 led to an era of intense research both in the UK and the US, known as the 18 first golden age of aeroacoustics [12]. An important result that came out of 19 this era is that jet noise scales with the eight power of the jet velocity [13, 14], 20 whereas the jet propulsive power only varies with the third power of the jet 21 velocity. Accordingly, it is possible to reduce the jet noise by lowering the jet 22 velocity and increasing the diameter of the jet. This led to the introduction 23 of high-bypass ratio turbofan engines, which today are in widespread use in 24 commercial aviation. For a comprehensive overview of the theory developed 25 during the first golden age of aeroacoustics, the interested reader is referred 26 to [15]. 27

As computers became more and more powerful, numerical predictions 28 of aerodynamically generated noise started to become possible. An early 29 example of this is the propeller and rotorcraft prediction codes developed 30 at NASA Langley [16, 17]. These codes did not explicitly resolve the flow 31 field, but instead required input from correlations, simplified aerodynamic 32 calculations, or experiments in order to estimate the loading on the blades, 33 which in turn translates into noise sources in the Ffowcs Williams - Hawkings 34 equation [18]. Later, in 1992, Sir Michael James Lighthill foresaw a second 35 golden age of aeroacoustics, this time driven by direct computations of the 36 aerodynamically generated noise sources using CFD [12]. This led to a new 37 field of aeroacoustics, known as computational aeroacoustics (CAA). The 38 progress of this field over the past three decades has been summarized in 39 several review papers, see e.g. [19, 20] 40

In general, CAA shares a lot of similarities with CFD, including the

need for robust numerical methods and accurate turbulence models for high-42 Reynolds number flows. However, as pointed out by e.g. Tam [21], some of 43 the challenges faced by CAA are often not encountered in CFD, including 44 the need to accurately resolve the acoustic waves (whose amplitude often are 45 orders of magnitude smaller than the amplitude of the hydrodynamic waves), 46 the need to propagate the acoustic waves over very large distances without 47 significant dissipation and dispersion errors, and the need to resolve unsteady 48 phenomena over a wide frequency range. In addition to this, reflections 49 of acoustic or hydrodynamic waves against an artificial boundary, e.g. an 50 inlet or an outlet to the computational domain, can quickly contaminate the 51 acoustic part of the solution, rendering the whole simulation useless [22]. 52 Therefore, numerical methods that are tailored to the needs of CAA have 53 been actively developed over the past decades. These include, but are not 54 limited to, the dispersion relation preserving (DRP) finite difference schemes 55 by Tam and Webb [23], the compact finite difference schemes by Lele [24], 56 the prefactored compact finite difference schemes by Ashcroft and Zhang 57 [25], the Navier-Stokes characteristic boundary conditions by Poinsot and 58 Lele [26], the radiation boundary conditions by Tam and Dong [27], Dong 50 [28], and the non-reflecting boundary conditions by Giles [29]. 60

The spatial discretization schemes mentioned above were all developed for structured grids, and some of them rely on wide finite-difference stencils. Although highly successful, the limitation of these schemes to structured grids makes it hard to adopt them for more complex configurations. Unstructured finite-volume schemes, on the other hand, are typically limited to secondorder accuracy in space. This raises the need for high-order unstructured schemes in order to tackle complex configurations of industrial interest with
optimal efficiency [20].

The development of high-order methods for unstructured grids has seen 69 great progress over the past decades. Interestingly, much of this development 70 has not been done with aeroacoustics in mind, but rather for the purpose 71 of performing accurate direct numerical simulations (DNS) and large eddy 72 simulations (LES) of turbulent flows. Among the many high-order schemes 73 that have been developed, we find the discontinuous Galerkin (DG) method 74 [30, 31], the flux reconstruction (FR) method [32], the spectral-difference 75 (SD) method [33, 34, 35, 36], and the continuous Galerkin (CG) method 76 [37].77

In recent years, the aforementioned high-order methods have started to be applied for CAA, see, e.g., [38, 39, 40, 41, 42]. However, given the large number of possible applications for CAA mentioned earlier, the potential benefits of these methods remain largely unexplored. Therefore, there is a need to further develop and test these methods in the context of CAA.

The purpose of this paper is to document a recently developed com-83 putational framework for jet noise predictions based on the high-order DG 84 method [40, 42]. This framework is based on a hybrid approach, in which 85 the turbulent jet is first simulated using the open-source spectral/hp element 86 framework Nektar++ [43, 44]. After this, the far-field noise is computed us-87 ing the Antares library [45, 46, 47]. Antares solves Formulation 1A [48] of 88 the Ffowcs Williams and Hawkings [18] equation in the time-domain using a 89 source-time dominant algorithm. 90

91

The present paper focuses on describing the setup of the framework using

Nektar++ and Antares. In particular, in section 2, the pre-processing steps (section 2.1), numerical method (section 2.2), and the post-processing steps (section 2.3) are described in detail. In section 4, the framework is then applied to an isolated and installed jet (section 4.1), followed by a demonstration of the scaling properties of the compressible flow solver in Nektar++ (section 4.2). Finally, in section 5, some conclusions are drawn.

98 2. Methodology

99 2.1. Pre-Processing

In this section we will present the pre-processing steps that we use to set up the jet noise simulations. An overview of these steps, including how they are linked together, is provided in Fig. 1.

103 2.1.1. CAD Preparation

Similar to many other computer aided engineering (CAE) problems, the 104 starting point for our simulations is a CAD description of the geometry that 105 we want to analyze. Typically, this geometry contains some surfaces that 106 are not relevant to the aeroacoustic analysis, such as small features that we 107 do not wish to resolve and/or surfaces that are not in direct contact with 108 the fluid. In addition to this, the outer boundaries of the computational 109 domain are often missing from the CAD. Therefore, the first step in the pre-110 processing pipeline is to clean up the CAD and add additional surfaces that 111 define the missing boundaries. To this end, we use the open-source software 112 Gmsh [49], which provides access to the OpenCASCADE CAD kernel. The 113 Gmsh software is in turn controlled through its Python API, which makes it 114 very convenient to automate the CAD preparation steps. 115



Figure 1: Flowchart illustrating the pre-processing setup.

116 2.1.2. Mesh Generation

The next step in the pre-processing pipeline is to generate a linear volume mesh. To this end, we use the commercial software STAR-CCM+ v2019.2. In particular, we use the *Advancing Layer Mesher* and the *Tetrahedral Mesher* in STAR-CCM+ to generate a single prism layer along solid walls, and tetrahedral elements in the rest of the domain, respectively.

The use of an unstructured grid has two important advantages. Firstly, it can be used to mesh complex geometries with relative ease, which is important in industrial applications. Secondly, it is straightforward to locally refine the mesh where it is needed, without having to refine other parts of the domain due to the constraints imposed by a structured grid. On the other

hand, when unstructured grids are used, it is very difficult to create high-127 quality anisotropic mesh elements, which could be beneficial in some areas, 128 such as the early shear layer of the jet. However, recent work has found that 129 it is beneficial to use isotropic mesh elements in the early shear layer for jet 130 noise applications [50, 51]. Therefore, we believe that the current setup is 131 well suited for our purposes. If, however, we would like to switch to a hybrid 132 meshing strategy in the future, it is possible since Nektar++ also supports 133 hexahedral and pyramid elements. 134

After the linear volume mesh has been generated, the NekMesh utility [44], which is part of the Nektar++ framework, is used to generate a highorder mesh with several prism elements in the wall-normal direction. To this end, the mesh is first projected onto the CAD by curving the mesh faces that are in contact with the boundary such that they conform to the CAD surface. After this, the curved prism elements are split in the wall-normal direction using an isoparametric approach [52].

If the mesh elements attached to the boundary are highly anisotropic 142 and/or the underlying geometry is highly curved, it is possible that self-143 intersecting elements are created during the CAD-projection step. To mini-144 mize the risk of this happening, the aspect ratio of mesh elements adjacent 145 to a boundary must be carefully controlled during the meshing process. In 146 addition to this, the prism elements should be curved before they are split in 147 the wall-normal direction. For complex geometries, however, it is often hard 148 to completely eliminate the presence of invalid elements. To ensure that the 149 simulations can still be run, the NekMesh utility therefore provides a mod-150 ule which can detect and remove the curvature from invalid elements. For 151

details on this, the interested reader is referred to the official documentation[53].

The output of the mesh projection step is a high-order mesh that conforms 154 to the underlying CAD geometry. This mesh is written to an HDF5 file [54] 155 using the layout described in [44]. The HDF5 format is used since this 156 enables several processes to read parts of the mesh in parallel, which in turn 157 is necessary to efficiently partition the mesh in massively parallel simulations. 158 We finally note that the NekMesh utility can read mesh files in both the 159 .ccm format used by STAR-CCM+ and the .msh format used by Gmsh. 160 Therefore, any meshing software that supports one of these file formats can 161 be used in the pre-processing pipeline shown in Fig. 1. 162

163 2.1.3. Initial Conditions

In this work, we use the commercial software STAR-CCM+ v2019.2 to 164 compute the initial condition for the LES simulations. More precisely, STAR-165 CCM+ is used to solve the Reynolds Averaged Navier Stokes (RANS) equa-166 tions using the $k - \omega$ SST turbulence model. For the RANS simulations, 167 we use the same mesh topology as for the LES simulations. However, since 168 STAR-CCM+ uses a finite volume discretization, whereas Nektar++ uses 169 a high-order discontinuous Galerkin discretization, we typically use a finer 170 mesh for the RANS simulations. 171

In order to interpolate the RANS solution onto the high-order expansion used in the LES simulations, the RANS solution is first interpolated to the mesh vertices and written to a .csv file. After this, we use the FieldConvert utility, which is part of the Nektar++ framework, to interpolate the point cloud defined by the .csv file onto the polynomial expan-

sion used in the LES simulations. More precisely, the field values are first 177 interpolated onto the quadrature points inside each element, followed by an 178 elemental projection in which the point values are converted to the corre-179 sponding modal coefficients. Since the RANS solution is relatively smooth, 180 and because it only represents the initial conditions for the LES simulations, 181 we typically project it onto a linear polynomial basis (p = 1). More details 182 about how to use the interpolation routines in FieldConvert can be found 183 in the official documentation [53]. 184

The interpolated solution is then written to a HDF5 file [54] using the layout described in [44]. The HDF5 format is used to export the high-order solution for the same reason that it is used to export the mesh, namely to enable parallel I/O where each process only reads a subset of the solution during initialization of the solver.

It is important to emphasize that there are other options for defining 190 initial conditions to the simulation. In particular, if another RANS solver 191 is available, it can easily be plugged into the pre-processing pipeline shown 192 in Fig. 1 as long as it can export the solution in the .csv format. In 193 addition to this, Nektar++ allows the user to define the initial conditions 194 using analytical expressions. Note that, in this case, there is no need to 195 write the initial conditions to file. Instead, the initial conditions can be 196 added directly to the input files for Nektar++. For more details on this, the 197 interested reader is referred to the official documentation [53]. 198

199 2.2. Numerical Method

The flow is modeled by the compressible Navier-Stokes equations written in conservative form

$$\frac{\partial \boldsymbol{q}}{\partial t} + \frac{\partial}{\partial x_i} (\boldsymbol{f}_i^c(\boldsymbol{q}) - \boldsymbol{f}_i^v(\boldsymbol{q}, \boldsymbol{\nabla} \boldsymbol{q})) = 0, \qquad (1)$$

where $\boldsymbol{q} = (\rho, \rho u_1, \rho u_2, \rho u_3, E)^T$ is the vector of conservative variables, $\boldsymbol{f}_i^c(\boldsymbol{q})$ is the convective flux in the i^{th} direction and $\boldsymbol{f}_i^v(\boldsymbol{q}, \nabla \boldsymbol{q})$ is the diffusive flux in the i^{th} direction. In this work, we assume that the gas is calorically perfect and obeys the ideal law. In addition to this, we assume that the viscosity is constant and equal to the far-field value outside the jet. This assumption is valid since we only consider cold jets whose temperature is close to the ambient temperature.

207 2.2.1. Spatial Discretization

The DG method is used to discretize Eq. (1) on a fully unstructured mesh [55]. The diffusion terms are treated with the incomplete Interior Penalty (IP) method [56] and the advection terms are rewritten using a standard weak DG scheme, followed by the application of Roe's approximate Riemann solver to compute the convective flux across the element boundaries [57, 58]. A detailed explanation of how to define the spatial discretization in Nektar++ is provided in the official documentation [53].

In traditional LES, subgrid-scale models are used to model the coupling between the resolved and the unresolved turbulent scales. However, by using the Roe Riemann solver for coupling between adjacent elements, the solution is inherently diffused by the numerical discretization introduced by the upwinding. This dissipation tends to smooth out high-frequency features and suppress small-scale turbulent fluctuations. As a result, the unresolved turbulent scales are implicitly handled through the dissipation introduced by the discretization scheme. Previous work has shown that the DG discretization in combination with Roe's approximate Riemann solver is well suited for implicit Large Eddy Simulations (iLES) [59, 60, 61, 62, 63, 64, 65].

In order to stabilize the solution in regions where the resolution is not high enough, an artificial viscosity approach is used. This stabilization strategy works by augmenting the dynamic viscosity and the thermal conductivity by an artificial viscosity and thermal conductivity in under-resolved regions. The artificial viscosity and thermal conductivity are respectively defined as

$$\mu_{av} = \mu_0 \rho \frac{h}{P} (c + \sqrt{u_k u_k}) S, \qquad (2)$$

$$\kappa_{av} = \mu_{av} \frac{C_p}{Pr},\tag{3}$$

where μ_0 controls the magnitude of μ_{av} , S is the sensor, C_p is the specific heat and Pr is the Prandtl number.

A sensor is used to make sure that the artificial viscosity and thermal conductivity are only activated in under-resolved regions. The sensor used in this work is a modal resolution based sensor [66]. To compute this sensor, the following resolution indicator is first computed.

$$s_e = \log_{10} \left(\frac{\langle u - \hat{u}, u - \hat{u} \rangle}{\langle u, u \rangle} \right), \tag{4}$$

where $\langle \cdot, \cdot \rangle$ is the L_2 inner product, u is the full expansion of a state variable (density in our case) of order P. Lastly, \hat{u} is the truncated expansion with terms up to P-1. Based on the resolution indicator, the sensor is computed as

$$S = \begin{cases} 0 & s_e < s_0 - \kappa, \\ \frac{1}{2} \left(1 + \sin\left(\frac{\pi(s_e - s_0)}{2\kappa}\right) \right) & |s_e - s_0| \le \kappa, \\ 1 & s_e > s_0 - \kappa, \end{cases}$$
(5)

where $s_0 = s_k - 4.25 \log_{10}(P)$. Note that, the parameters s_{κ} and κ have to be chosen in the correct range in order to make the artificial viscosity effective.

229 2.2.2. Temporal Discretization

The set of ordinary differential equations (ODEs) obtained from the spa-230 tial discretization are integrated in time using a second-order singly diago-231 nally implicit multi-stage Runge–Kutta (SDIRK2) method [67]. The nonlin-232 ear system from SDIRK2 is iteratively solved by the Jacobian-free Newton 233 Krylov (JFNK) method [55]. Thus, the Newton method is used to solve the 234 nonlinear system [68]. The restarted generalized minimal residual method 235 (GMRES) [69] is adopted to solve the linear system. A Jacobian-free method 236 is applied in order to avoid explicitly calculating and storing the Jacobian 237 matrix [68]. 238

In addition, a preconditioner in GMRES is employed which is important 239 for solving stiff problems. The parameters used to set up the preconditioner 240 are shown in listing 1, where PreconMatFreezNumb specifies number of 241 iteration that the GMRES solver will perform before the preconditioner ma-242 trix be updated, NonlinIterTolRelativeL2 is the convergence toler-243 ance of the nonlinear system relative to the initial nonlinear system residual, 244 LinSysRelativeTolInNonlin represents the convergence tolerance of 245 the linear system solver in each nonlinear iteration. PreconItsStep de-246

termines the number of preconditioning iterations to calculate the preconditioned vector. For further understanding of the mathematical formulation of
these parameters, the reader may find useful the work of Yan et al. [55].

It is worth mentioning that the parameters shown in Listing 1 have a 250 strong impact on the convergence and efficiency of the implicit solver, and 251 consequently speed up the simulation. Special attention should be given 252 to the PreconMatFreezNumb parameter, since if its value is too low the 253 preconditioner matrix may be updated in every time-step, which reduces 254 the efficiency. However, large values may lead to divergence since the pre-255 conditioner matrix is not updated as frequently as the numerical simulation 256 requires, making the preconditioner less effective at every time step. The 257 PreconItsStep parameter may significantly reduce the number of GM-258 RES and nonlinear iterations when its value is increased; the time per time-259 step increases but it allows for larger time steps which speeds up the simu-260 lation overall. Therefore, the appropriate balance of these two parameters is 261 essential for the efficient execution of the numerical simulation. 262

```
<SOLVERINFO>
263
        <!-- This part is put inside the <CONDITIONS> tag -->
264
        PreconMatFreezNumb
                                        = 1000
                                                   </P>
265
                                                   </P>
        NonlinIterTolRelativeL2
                                        = 1.0E - 3
266
        <P> LinSysRelativeTolInNonlin = 5.0E-2
                                                   </P>
267
        PreconItsStep
                                                  </P>
                                        = 3
268
    </SOLVERINFO>
269
```

Listing 1: Parameters for the preconditioner.

270 2.2.3. Boundary Conditions

Boundary conditions can have a large impact on the stability and accuracy of a simulation. In LES simulations of jet noise, three types of boundaries are typically present; solid walls, far-field boundaries, and inlet(s) to the
nozzle. In this section we will describe the implementation of these boundary
conditions in detail.

The far-field boundaries are used to impose the ambient conditions of the 276 flow, far away from the jet. The way the ambient conditions are imposed 277 depends on whether the flow enters or exits the domain though the boundary. 278 Along the parts of the far-field boundary where the flow enters the domain, 279 the far-field state is imposed using the Weak-Riemann approach described 280 in [57]. In this approach, the inviscid flux across boundary faces is computed 281 from the solution of a Riemann problem, using the far-field state and the 282 internal solution as the "left" and "right" state, respectively. This boundary 283 condition is non-reflective for normally incident waves as long as the Rie-284 mann solver correctly differentiates between incoming and outgoing waves. 285 The Riemann solver used in this work approximately satisfies this condition. 286 Despite not being perfectly non-reflective, the amount of reflections against 287 this boundary is deemed to be negligible since the far-field boundary is placed 288 far away from the jet, and because the mesh is coarsened as it approaches 280 the far-field boundary. 290

At the outlet part of the far-field boundary, the *Weak-Riemann* approach is also used. However, in contrast to the inlet part of the boundary, only the ambient pressure is imposed. This is done by setting the left state in the Riemann problem to be the internal solution, but with the pressure substituted for the ambient pressure [57]. This boundary condition is reflective. Therefore, it is important that all waves are attenuated before they reach this boundary. If only mesh coarsening is used for this purpose, an excessively long computational domain would need to be used. To avoid this, we add a
sponge zone upstream of the outlet [70]. In this zone, the right hand side of
Eq. (1) is augmented with a damping term given by

$$\dots = -\sigma(\boldsymbol{x}) \left(\boldsymbol{q} - \overline{\boldsymbol{q}}\right). \tag{6}$$

Here, $\sigma(\boldsymbol{x})$ is a function that controls the magnitude of the damping and $\overline{\boldsymbol{q}}$ is a steady reference solution. In this work, the RANS solution that we use to initialize the simulations is also used to define $\overline{\boldsymbol{q}}$.

As explained in [22], the best results are obtained when the function $\sigma(\boldsymbol{x})$ is increased slowly over a sufficiently long distance. To achieve this, we compute it as

$$\sigma(\boldsymbol{x}) = \begin{cases} 0 & x_1 < x_{s_1}, \\ \sigma_{\max} S\left(\frac{x_1 - x_{s_1}}{x_{s_2} - x_{s_1}}\right) & x_1 \in (x_{s_1}, x_{s_2}), \\ \sigma_{\max} & x_1 > x_{s_2}. \end{cases}$$
(7)

Here, (x_{s_1}, x_{s_2}) denotes the interval where the sponge is increased (the jet is aligned with the x_1 -axis), σ_{\max} is the maximum value of the damping coefficient, and $S(\xi)$ is the smooth-step function

$$S(\xi) = 6\xi^5 - 15\xi^4 + 10\xi^3, \quad \xi \in (0, 1).$$
(8)

The smooth-step function is shown in Fig. 2. As can be seen from this figure, it increases smoothly from 0 to 1 over the interval (0, 1).

The sponge zone is activated in Nektar++ by adding the commands shown in Listing 2 and Listing 3 under the <CONDITIONS> tag and the



Figure 2: Smooth-step function.

314 <FORCING> tag in the input file, respectively. Note that these commands 315 require that the parameters x1, x2, and sigma_max have been defined under 316 the <PARAMETERS> tag in the input file.

```
<!-- This part is put inside the <CONDITIONS> tag -->
317
    <FUNCTION NAME="AverageSolution">
318
        <F VAR="rho, rhou, rhov, rhow, E" FILE="avg.fld" />
319
    </FUNCTION>
320
321
    <FUNCTION NAME="DampingCoeff">
322
        <E VAR="rho,rhou,rhov,rhow,E" VALUE="-sigma_max*((x>x1)*(x<x2)*(6*((x-x1)/(x2</pre>
323
         -x1))^5 - 15*((x-x1)/(x2-x1))^4 + 10*((x-x1)/(x2-x1))^3) + (x>=x2))" />
324
    </FUNCTION>
325
```

Listing 2: Definition of average solution and damping coefficient.

```
326 <!-- This part is put inside the <FORCING> tag -->
327 <FORCE TYPE="Absorption">
328 <REFFLOW> AverageSolution </REFFLOW>
329 <COEFF> DampingCoeff </COEFF>
330 </FORCE>
```

Listing 3: Definition of sponge zone.

The length of the sponge zone is set to $20D_j$, where D_j is the diameter of the jet. The damping coefficient, which has a unit equal to s⁻¹, is then selected such that it is proportional to the reciprocal of the residence time of a convected disturbance inside the sponge zone. For the simulations presented in this work, we found that a value of

$$\sigma_{\max} = 10 \frac{U_j}{x_{s_2} - x_{s_1}},\tag{9}$$

where U_j is the velocity of the jet at the nozzle exit, worked well.

At the inlet to the nozzle, we impose the stagnation pressure, stagnation temperature, and the flow direction. This boundary condition is chosen because in most experiments, including the ones used to validate the simulations performed in this work, the stagnation conditions are known upstream of the nozzle exit. The flow direction, on the other hand, is typically not known. Therefore, we set the flow direction equal to the boundary normal direction.

The inlet boundary condition is also imposed weakly through the Rie-344 mann solver. To this end, the 4 variables imposed by the boundary condition 345 must be combined with one variable from the interior in order to form a com-346 plete left state for the Riemann solver. Depending on which interior variable 347 is "extrapolated", the flow will develop in a different way from a given initial 348 condition. For the simulations presented in this paper, it was found that 349 extrapolating the magnitude of the velocity gave good stability. More pre-350 cisely, the left state in the Riemann solver is computed as follows. First, the 351 temperature at the boundary is computed from the imposed stagnation tem-352 perature and the magnitude of the velocity taken from the interior, following 353

³⁵⁴ the definition of the stagnation temperature

$$T_{\rm BC} = T_{0,\rm BC} - \frac{1}{2C_p} u_{\rm R}^2.$$
 (10)

Here, $u_{\rm R} = \sqrt{u^2 + v^2 + w^2}$ is the magnitude of the velocity obtained from the interior solution and C_p is the specific heat capacity. After this, the density is computed using the well known isentropic flow relation

$$\rho_{\rm BC} = \rho_{0,\rm BC} \left(\frac{T_{\rm BC}}{T_{0,\rm BC}}\right)^{\frac{1}{\gamma-1}},\tag{11}$$

where R is the specific gas constant, γ the ratio of specific heats, and $\rho_{0,BC} = \frac{p_{0,BC}}{RT_{0,BC}}$ is the stagnation density. Finally, the left state at the boundary is defined as

$$\boldsymbol{q}_{\mathrm{L}} = \begin{bmatrix} \rho_{\mathrm{BC}} \\ \rho_{\mathrm{BC}} u_{\mathrm{R}} \boldsymbol{n} \\ \frac{p_{\mathrm{BC}}}{\gamma - 1} + \frac{1}{2} \rho_{\mathrm{BC}} u_{\mathrm{R}}^2 \end{bmatrix}.$$
 (12)

Here, $\boldsymbol{n} = (n_1, n_2, n_3)$ is the imposed flow direction and $p_{\rm BC} = \rho_{\rm BC} R T_{\rm BC}$ is the static pressure.

To use the inflow boundary condition described above, the commands shown in Listing 4 are added under the <BOUNDARYCONDITIONS> tag in the input file. In this listing, rho0 is the stagnation density, (n1, n2, n3) define the flow direction, and E0 = $p_{0,BC}/(\gamma - 1)$ is the stagnation total energy. These parameters must be defined inside the <PARAMETERS> tag in the input file.

^{369 &}lt;!-- This part is put inside the <BOUNDARYCONDITIONS> tag -->
370 <REGION REF="ID">
371 <D VAR="rho" USERDEFINEDTYPE="StagnationInflow" VALUE="rho0" />

```
372 <D VAR="rhou" USERDEFINEDTYPE="StagnationInflow" VALUE="n1" />
373 <D VAR="rhov" USERDEFINEDTYPE="StagnationInflow" VALUE="n2" />
374 <D VAR="rhow" USERDEFINEDTYPE="StagnationInflow" VALUE="n3" />
375 <D VAR="E" USERDEFINEDTYPE="StagnationInflow" VALUE="n3" />
376 </REGION>
```

Listing 4: Definition of stagnation inflow boundary condition.

To save computational resources, we only include a short part of the nozzle 377 in our simulations. As a result, the boundary layer does not have enough 378 time to develop between the inlet and the end of the nozzle. To overcome 379 this issue, we impose non-uniform profiles of stagnation temperature and 380 pressure at the inlet. This can be done by passing a file containing the 381 boundary expansion of ρ_0 and E_0 to the boundary condition. Alternatively, 382 analytical expressions describing the spatial variation of ρ_0 and E_0 can be 383 used instead of the constant values shown in Listing 4. 384

To construct a file containing the boundary expansion of ρ_0 and E_0 , we start by extracting the mesh elements defining the boundary using the extract module in NekMesh (see section 4 in the official documentation for details [53]). After this, the resulting mesh file is edited to ensure that the boundary expansion defined in the file uses the same polynomial order as the one that will be used in the simulations. An example of this is for a mesh containing quad and triangular elements is provided in Listing 5.

```
392 <!-- This part is put inside the <NEKTAR> tag -->
393 <EXPANSIONS>
394 <E COMPOSITE="C[0]" TYPE="MODIFIED" NUMMODES="2" FIELDS="rho,E" />
395 <E COMPOSITE="C[1]" TYPE="MODIFIED" NUMMODES="2" FIELDS="rho,E" />
396 </EXPANSIONS>
```

Listing 5: Expansion definition for an inlet containing quad and triangular elements.

After the expansion definition has been edited, the quadrature points are 397 exported to a .csv file using FieldConvert. By default, FieldConvert 398 will export data on a uniform set of points for visualization purposes. How-399 ever, since we want to project data on the quadrature points back onto 400 the polynomial expansion, we need to export the location of the quadrature 401 points instead. To do this, the flag --noequispaced should be passed 402 to FieldConvert. After this, we use Python to compute the values of ρ_0 403 and E_0 at the location of the quadrature points. Note that, if an analytical 404 expression is used for this purpose, it is simpler to specify this expression 405 directly in the input file. In this work, however, we specify the profiles of 406 ρ_0 and E_0 using experimental data. More precisely, we use SciPy [71] to 407 interpolate the experimental data to the location of the quadrature points. 408 After this, we write the coordinates of the quadrature points together with 409 the associated values of ρ_0 and E_0 to a new .csv file. A new field file con-410 taining the boundary expansions of ρ_0 and E_0 can then be obtained using 411 the pointdatatofld module in FieldConvert. 412

The implementation of the solid wall-boundary condition used for the nozzle walls is described in detail in [57].

415 2.3. Post-Processing

In this section we describe the different tools that we use to post process the solution. In particular, the technique used to propagate the noise between the jet and the far-field observers is described in section 2.3.2. It does in turn depend on the data collection strategy described in section 2.3.1.

420 2.3.1. Data Collection

Nektar++ provides a set of filters that can be used to compute derived 421 quantities from the solution. In this work, we make extensive use of the 422 history points filter to probe the solution at a set of predefined points. By 423 default, the history point filter writes the data to a .csv file every N^{th} time 424 step. If a small number of points is used, this is usually not a problem. 425 However, as more and more points are added, the I/O quickly starts to 426 consume a significant portion of the overall execution time. In addition to 427 this, the .csv format can not be used to store complex data structures, it 428 requires a relatively large amount of memory per data point, and it does not 429 allow the user to easily add additional metadata to the file. To overcome 430 these limitations, the history point filter in Nektar++ was extended to store 431 the points in the HDF5 format [54]. 432

The HDF5 format is a flexible data format that allows the user to design the layout of the file based on the needs of a particular application. To this end, the HDF5 group provides the HDF5 library, which gives the user access to a rich API that can be used to create, structure, and read/write large amounts of data to the file.

The file structure chosen to store the history points is illustrated in Fig. 3. The folders shown in this figure represent HDF5 groups. These groups are synonymous to folders in a file system, in the sense that they can be used to organize data inside the file. Starting from the root group, which we call NEKTAR, two sub-groups are added. The first is called COORDINATES which, as the name suggests, contains the coordinates of history points stored as three HDF5 datasets. The second sub-group below the root group is called

TIME-DATA. Each time a new time step is written to the file, a new sub-445 group called TIME-STEP-N is added to the TIME-DATA group, where N 446 is an integer which is incremented by one each time a new time step is 447 added. Inside this sub group, five HDF5 datasets are created, one per solution 448 variable. In addition to this, the current time is added as a HDF5 attribute. 449 The HDF5 format is well suited for post-processing since many popular 450 high-level programming languages, such as Python, Julia, and MATLAB, can 451 be used to read/write HDF5 files. Support for this functionality is usually 452 provided either directly, or through high-quality third-party packages. In 453 this work, we have used the h5py package to post-process the history point 454 data in Python. One example of this is provided in the next section. 455

456 2.3.2. Far-Field Noise Predictions

The overall goal of this work is to predict the far-field noise of turbulent jets. Usually, the location where we want to measure the noise is located outside the computational domain. Therefore, a separate method is needed to propagate the noise between the jet and the far-field. In this work, we use Antares 1.17.0 [45, 46, 47] for this purpose.

Antares solves the Ffowcs Williams - Hawkings equation [18] in the time 462 domain using a source-time dominant algorithm. The solution to the Ffowcs 463 Williams - Hawkings equation expresses the pressure signal at an arbitrary 464 observer location as an integral over a closed surface, plus a volume integral 465 over the volume outside the surface. In this work, we don't include the 466 volume integral in the solution to the Ffowcs Williams - Hawkings equation. 467 This simplification is widely used in jet noise applications [10, 11, 72], and 468 is valid as long as the integration surface is chosen such that it encloses 469



Figure 3: Layout of an HDF5 file storing history point data.

all relevant noise sources [73], and such that no entropy or vorticity waves 470 cross the surface [74, 75]. To satisfy these conditions, a conical surface that 471 follows the spreading rate of the jet is usually chosen [11]. Ideally, this 472 surface should be placed as close to the jet as possible in order to minimize 473 numerical dissipation and dispersion of the acoustic waves as they propagate 474 between the source and the surface. However, if the surface is too narrow, 475 the conditions stated above will be violated. Therefore, there is an optimal 476 width of the integration surface [10, 11]. 477

In addition to adjusting the width, the length of the surface will influence 478 the accuracy of the results. Typically, a length of 30 nozzle diameters is 479 considered enough to enclose all relevant noise sources [72]. Unfortunately, 480 the entropy and vorticity waves generated in the jet do not decay over this 481 distance. Therefore, if the surface is closed at the downstream end, the sec-482 ond condition stated above will be violated. As a result, spurious noise is 483 generated by the "end-cap" of the surface, which degrades the quality of the 484 results. One way to solve this problem is to extend the domain and the 485 integration surface much further, but since this also increases the compu-486 tational cost, it is not desirable. Instead, alternative strategies such as the 487 method of end-caps [74], the pressure-formulation of the Ffowcs Williams 488 - Hawkings equation [74, 75], and the additional surface terms developed 489 by Rahier et al. [76], have been proposed. Unfortunately, the method of 490 end-caps and the additional surface terms proposed by Rahier et al. [76] are 491 not supported by Antares. In addition to this, the pressure formulation was 492 found to not improve the results for the cold jets considered in this work [40]. 493 Therefore, a fourth strategy, in which the integration surface is left open at 494

the downstream end, is adopted instead. The benefit of this method is that it eliminates spurious noise from the end-cap. However, this comes at the cost of less accurate noise predictions for the lowest frequencies, see [74] for details.

At present, Antares provides two solutions to the Ffowcs Williams -Hawkings equation, Formulation 1A by Farassat [48] and Formulation 1C by Najafi-Yazdi et al. [77]. The main difference between these two formulations is that the latter includes the effect of a flight stream on noise propagation. In this work, we only consider jets operating in a medium at rest. Therefore, Formulation 1A is used.

To compute the noise propagation, we need the time history of the solution on the integration surface. The process that we use to generate this time history is illustrated in Fig. 4. It is important to point out that this process is relatively generic, and will therefore work with any post-processing software that implements the Ffwocs Williams - Hawkings method, as long as this software can read one of the file formats that Nektar++ exports.

To begin with, the integration surface is constructed in Gmsh [49] using 511 the OpenCASCADE CAD kernel. After this, Gmsh is also used to create a 512 triangular surface mesh. The mesh resolution on the integration surface is 513 typically higher than the volume mesh. The reason for this is that Antares 514 does not support Gaussian quadrature on high-order elements. Instead, it 515 stores the data on the mesh vertices, and then compute the element con-516 tribution as the average over all vertex values, scaled by the area of the 517 element. 518

⁵¹⁹ To sample data on the integration surface, we use the history point filter

described in section 2.3.1. We start by converting the mesh generated with Gmsh into the format used by Nektar++ using the NekMesh utility. By default, NekMesh adds an XML tag to the mesh file that defines which type of polynomial expansion, and which quadrature rule, to use for each composite (group of mesh elements). Since Antares does not support highorder elements, we edit the default expansion definition such that linear finite elements are used. An example of this is provided in listing 6.

```
527 <!-- This part is put inside the <NEKTAR> tag -->
528 <EXPANSIONS>
529 <E COMPOSITE="C[0]" TYPE="MODIFIED" NUMMODES="2" FIELDS="u" />
530 </EXPANSIONS>
```

Listing 6: Expansion definition for the Ffowcs Williams - Hawkings integration surface.

After the expansion definition has been edited, the quadrature points are 531 exported to a .csv file using the FieldConvert utility. As explained in 532 section 2.2.3, FieldConvert exports data on a uniform set of points instead 533 of the quadrature points by default. For a linear finite element expansion, 534 these points are simply the vertices of the elements. Although this is the data 535 that we eventually want to export to Antares, we do not want to sample the 536 solution at these points for reasons that will become clear soon. Therefore, 537 the flag --noequispaced is passed to FieldConvert when the points 538 are exported to the .csv file. The points defined in the .csv file are then 539 used by the history point filter to sample the solution at the integration 540 surface. 541

Once the simulation is finished, we need to convert the history point data into a format that Antares can read. To this end, the Python script shown in listing 7 is used. This script works by looping through all time steps stored

in the HDF5 file using the h5py package. For each time step, the script 545 converts the solution to primitive variables, and stores the result in a tem-546 porary .csv file. The .csv file is then used by the process module called 547 pointdatatofld, which projects data on the Gaussian quadrature points 548 onto an existing expansion. Here, the existing expansion is stored in the 549 field variable in the script, which in turn is obtained by reading the mesh 550 file described earlier. Once the projection is completed, the updated expan-551 sion, which now contains 5 fields (one for each primitive variable) is written 552 to a new file in the TecPlot binary format. Since the flag --noequispaced 553 is not passed to FieldConvert this time, the solution will be exported on 554 the mesh vertices instead of the quadrature points. Note that all these oper-555 ations are performed by directly calling the underlying Nektar++ functions 556 through the NekPy module, which is provided together with Nektar++. 557

```
from NekPy.FieldUtils import Field, InputModule, ProcessModule, OutputModule
558
    import h5py
559
560
    import numpy as np
561
562
    # Name of HDF5 file
    hdf5_filename = "fwh-data.hdf5"
563
564
    # Name of output file (in tecplot format)
565
566
    plt_filename = os.path.splitext("fwh-data.plt")
567
    # Name of mesh file (including expansion definition)
568
    mesh_filename = "fwh-surface.xml"
569
570
571
    # Open HDF5 file
    with h5py.File(hdf5_filename, 'r') as hdf5_file:
572
573
574
         # Read coordinates
        cgroup = hdf5_file["NEKTAR/COORDINATES"]
575
576
        x = cgroup["x"][()]
```

```
577
        y = cgroup["y"][()]
        z = cgroup["z"][()]
578
579
        # Go through all time steps
580
        tgroup = hdf5_file["NEKTAR/TIME-DATA"]
581
        for indx, tstep in enumerate(tgroup.values()):
582
583
584
             # Read solution at this time step
             rho = tstep['rho'][()]
585
             rhou = tstep['rhou'][()]
586
             rhov = tstep['rhov'][()]
587
             rhow = tstep['rhow'][()]
588
589
             Е
                  = tstep['E'][()]
590
             # Convert to primitive variables
591
             u = rhou / rho
592
             v = rhov / rho
593
             w = rhow / rho
594
             p = (E - 0.5 * (rhou **2 + rhov **2 + rhow **2)/rho) * 0.4
595
596
             # Save data in .csv format
597
             np.savetxt("tmp.csv", np.column_stack((x,y,z,rho,u,v,w,p)), delimiter=","
598
         , header="x, y, z, rho, u, v, w, p")
599
600
             # Create a Nektar++ Field and read expansion from file
601
             field = Field([])
602
603
             InputModule.Create("xml", field, mesh_filename).Run()
604
605
             # Project points to expansion
             ProcessModule.Create("pointdatatofld", field, frompts="tmp.csv").Run()
606
607
608
             # Save data to .plt format
             output_filename = plt_filename[0] + "-%05i"%(index) + plt_filename[1]
609
             OutputModule.Create("plt", field, output_filename).Run()
610
```

Listing 7: Python script for converting history point data into TecPlot binary format.



Figure 4: Flowchart illustrating the steps required to sample data on the integration surface used for the Ffowcs Williams - Hawkings method.

611 3. Performance considerations

Parallelization. Parallelization in Nektar++ is carried out using a domain 612 decomposition strategy, where the computational mesh is divided into one 613 subset per process. Mesh partitioning is achieved using either the Scotch 614 or the METIS libraries at the beginning of the simulation. Note that each 615 processor only reads the geometric information that is required for its own 616 partition, that is, the partitioning of the mesh file is done in parallel. Par-617 allel mesh partitioning, contrary to serial mesh partitioning, avoids one sin-618 gle thread reading the whole mesh and partitioning it alone which could 619 lead to the node running out of memory. This is achieved through the 620

use of HDF5 routines when reading the mesh datasets and the use of the --use-hdf5-node-comm flag with Nektar++. Output solution files can also be written in HDF5 format which is enabled by the flag --io-format Hdf5. The different processes communicate through the standard MPI protocol, where each process is independently executed and there is no explicit use of shared memory protocols such as OpenMP. In this work we use the Scotch library to partition the mesh.

Vectorization. The limitation in performance of modern hardware concerns 628 memory bandwidth speeds more than processor clock speeds. For this reason, 629 the use of arithmetically intense schemes, that is, schemes performing a high 630 number of floating point operations for each byte transferred from memory, 631 continues to be of considerable interest. At high polynomial orders, high-632 order or spectral/hp element methods are arithmetically intense schemes 633 since they involve dense, compact kernels for key finite element operators. 634 This is a significant advantage of high-order methods compared to lower-order 635 methods. Key aspects to exploit this potential for increased performance 636 are matrix-free implementations of finite element operators combined with 637 the use of sum-factorization and effective approaches for exploiting single-638 instruction multiple-data (SIMD) vectorization. Nektar++ is provided with 639 the matrix-free evaluation of basic finite element operations which still utilize 640 sum-factorization even for non-tensor-product elements (triangles, tetrahe-641 dra, and prisms) [78, 79, 80]. Nektar++ also allows explicit exploitation 642 of SIMD vectorization, showing that although performance is naturally de-643 graded when going from hexahedra to triangles, tetrahedra, and prismatic 644 elements, efficient implementations are still obtained, achieving 50% to 70%645

of the peak floating-point throughput of modern processors with both AVX2 646 and AVX512 instruction sets [81]. In addition to that, the SIMD vectoriza-647 tion of the diffusion operator and the block diagonal matrix obtained after 648 decoupling the jacobian matrix to calculate the preconditioned vector has 649 recently been implemented in Nektar++. Also, the $(\hat{L} + \hat{U})$ matrix, which 650 is also obtained from the jacobian matrix, employed to calculate the precon-651 ditioned vector is now able to use the auto-tuning routines (e.g. matrix free 652 operator) to optimize this operator. For further details about the precon-653 ditiner calculation the reader is referring to [55] 654

Automatic selection of the operations optimization. An auto-tuning capability, to automatically select the most efficient operations optimization for each execution, is also available in Nektar++ [80]. It can be turned on by including the lines, given in Listing 8, in the problem setting script.

```
660 <!-- This is how to enable auto tuning -->
661 <COLLECTIONS DEFAULT="auto"/>
662 </NEKTAR>
```

Listing 8: Flag for enabling performance auto-tuning in Nektar++.

For more details on this, the interested reader is referred to the official documentation [53].

665 4. Results

<NEKTAR>

659

666 4.1. Isolated and Installed Jet Noise

In this section we will present the application of the computational framework presented in this paper to investigating noise generated by isolated and installed jets.

670 4.1.1. Case Description

The nozzle considered in this work has an inner diameter of $D_j = 40$ mm, 671 a convergence half-angle of 2.44°, and a total length of $19D_j$. Both an isolated 672 and an installed configuration are considered. In the installed configuration, 673 an uniform-chord (2D) NACA4415 airfoil at 4° angle of attack is placed close 674 to the nozzle. The axial and vertical distance between the nozzle exit and 675 the airfoil trailing edge are $L = 3D_j$ and $H = 0.6D_j$, respectively. The chord 676 and span of the airfoil are $3.75D_j$ and $15D_j$, respectively. A schematic view 677 of the installed configuration is shown in Fig. 5. 678



Figure 5: Schematic view of the nozzle installed under the wing.

In this work, we consider a single operating point of the nozzle, corresponding to an acoustic Mach number of $M_a = U_j/c_{\infty} = 0.6$, a Reynolds number of $Re_j = \rho_{\infty}U_jD_j/\mu_{\infty} = 5.5 \cdot 10^5$, and a static temperature ratio of $T_j/T_{\infty} = 0.9335$. In addition to this, no flight stream is considered, meaning that the ambient medium is at rest.

Both the isolated and installed configuration of the nozzle have recently been tested in the Doak Laboratory Flight Jet Rig, located at the University of Southampton. During these tests, both aerodynamic data (velocity measurements in the jet plume) and far-field acoustic data was recorded. This data will be used in sections 4.1.3 and 4.1.4 to validate the simulations. More details about the experimental setup can be found in [40, 42, 82].

690 4.1.2. Computational Setup

A schematic view of the axi-symmetric, funnel-shaped computational do-691 main used in this work is presented in Fig. 6. Along the left and top boundary 692 shown in this figure, the far-field state is imposed using the *Weak-Riemann* 693 approached described in section 2.2.3. As indicated by the arrows in the 694 figure, a small co-flow corresponding to 2% of the jet exit velocity is also 695 added to the (stagnant) far-field state. This is done to ensure that vortical 696 structures generated in the jet are flushed out of the domain and to facilitate 697 flow entrainment. 698

⁶⁹⁹ Along the right boundary in Fig. 6, the ambient pressure is imposed ⁷⁰⁰ according to the method described in section 2.2.3. Since this boundary con-⁷⁰¹ dition is reflective, a sponge zone of length $20D_j$ is also added upstream of the ⁷⁰² outlet boundary. In this region, the damping term described in section 2.2.3 ⁷⁰³ is used to attenuate the jet before it reaches the outflow boundary.

The inlet to the nozzle is placed $1.5D_j$ upstream of the nozzle exit. At this boundary, non-uniform profiles of stagnation pressure and temperature are imposed to ensure that the mean velocity matches the experimental data as closely as possible at the nozzle exit, see section 2.2.3 for details about the implementation.

⁷⁰⁹ In addition to matching the mean velocity profile at the nozzle exit, re-⁷¹⁰ cent work has demonstrated the importance of getting the turbulence levels



Figure 6: Schematic view of the computational domain and the Ffowcs Williams - Hawkings integration surface.

in the boundary layer right [50, 51, 83, 84]. In order to satisfy this condi-711 tion, some type of artificial tripping mechanism and/or injection of synthetic 712 turbulence at the inlet boundary may be necessary when only a short part 713 of the nozzle is included in the simulations. In addition to this, for high 714 Reynolds number applications, it is usually necessary to model the inner 715 part of the boundary layer, either through the use of a wall model [50] or a 716 hybrid RANS-LES model [51, 85]. Unfortunately, Nektar++ currently does 717 not support any tripping mechanism, synthetic turbulence injection, or wall 718 model. Therefore, a standard no-slip boundary condition is used to model 719 the nozzle instead. In earlier work, we noted that the turbulence levels at 720 the nozzle exit are not accurately predicted with this setup [40, 42]. It is 721 well known that the state of the boundary at the nozzle exit can have a large 722 influence on the far-field acoustic results [50, 83, 84]. Therefore, a better wall 723 modeling strategy is expected to improve the results, but was unfortunately 724 outside the scope of this work. 725

The methodology described in section 2.1.2 was used to generate the two unstructured meshes shown in Fig 7. As can be seen from this figure, the main difference between these two meshes is that the latter includes the wing. In addition to this, the resolution far downstream is slightly lower for the second mesh. The resulting mesh count for the two meshes are summarized in Table 1. More details about the two meshes are provided in [42].

Three simulations have been performed, one for the isolated jet, and two for the installed jet. The simulation of the isolated jet was performed with Mesh-1 using a polynomial degree of P = 2. The two simulations of the installed jet were performed with Mesh-2, using a polynomial degree of



Figure 7: Illustration of unstructured grids defined in Table 1.

Name	Configuration	$N_{\rm prism}$	$N_{\rm tet}$
Mesh-1	Isolated	$4.36\cdot 10^5$	$6.18\cdot 10^6$
Mesh-2	Installed	$4.63\cdot 10^5$	$4.54\cdot 10^6$

Table 2: Simulation settings. Δt_{jet} and $\Delta t_{\text{FW-H}}$ denote the sampling rate in the jet plume and on the Ffowcs Williams - Hawkings integration surface, respectively. τ_{init} and τ_{sample} denote the time the simulation is run before sampling and during sampling, respectively.

Name	P	Mesh	DOF	Initialize from	$\frac{\Delta t c_{\infty}}{D_j}$	$\frac{\Delta t_{\rm jet} c_\infty}{D_j}$	$\frac{\Delta t_{\rm FW-H} c_{\infty}}{D_j}$	$\frac{\tau_{\text{init}}c_{\infty}}{D_j}$	$\frac{\tau_{\rm sample} c_\infty}{D_j}$
Isolated-1	2	$\operatorname{Mesh-1}$	$70\cdot 10^6$	RANS	0.002	0.04	0.04	350	320
Installed-1	2	Mesh-2	$54\cdot 10^6$	RANS	0.002	0.04	0.04	450	900
Installed-2	3	Mesh-2	$109\cdot 10^6$	Installed-1	0.002	0.04	0.04	300	450

P = 2 and P = 3, respectively. A summary of the settings used for each simulation is presented in Table 2.

738 4.1.3. Aerodynamic Results

We start by considering the mean and RMS of the axial velocity in the 739 jet plume. The simulation results obtained for the isolated jet are compared 740 against experimental data in Figs. 8a and 8b. As can be seen from these 741 figures, the agreement is quite satisfactory for all radial and axial locations. 742 The most notable discrepancies are found close to the jet centerline and fur-743 ther downstream. In this region, the mean velocity and turbulence levels are 744 over-predicted and under-predicted by the simulation, respectively. From 745 earlier work, we know that the state of the boundary layer at the nozzle exit 746 is not perfectly predicted in our simulations due to insufficient resolution and 747 lack of a wall-model [42]. Since the state of the boundary layer affects the 748 development of the shear layer, it is possible that the under-resolved bound-740 ary layer contributes to the discrepancies seen far downstream in Figs.. 8a 750 and 8b. Another possible source of the discrepancies is that we only include 751 a short part of the conical nozzle in the simulations, which in turn should 752

⁷⁵³ lead to slight discrepancies in the radial velocity component at the nozzle
⁷⁵⁴ exit. This is explained in more detail in [42].

The simulation results obtained for the installed nozzle are compared 755 against the corresponding experimental data in Fig. 8. In general, we can 756 note that the agreement between simulations and experiments is quite good 757 for the installed nozzle as well. It is also interesting to note that the higher 758 polynomial degree used in the simulation called "Installed-2" only has a mi-759 nor impact on the flow statistics. Finally, we note that the discrepancies be-760 tween the simulations and the experiments is consistent with those observed 761 for the isolated nozzle. This is expected, considering that the simulation 762 setup is almost identical for the two configurations. 763

764 4.1.4. Far-Field Acoustic Results

As explained in section 2.3.2, we use the Ffowcs Williams - Hawkings method implemented in Antares 1.17.0. to compute the far-field noise. The output of Antares is the pressure signal at the same microphone locations as were used in the experiments. The time-series signal is then converted into a power spectral density (PSD) in units [dB/St] using the implementation of Welch's method [86] in SciPy [71]. More details on the exact locations of the microphones and how the PSD is computed can be found in [40, 42].

The PSD at two representative microphone locations located at $\theta = 90^{\circ}$ and 43° with regards to the downstream jet axis is presented in Fig. 9a and 9b, respectively. By comparing the results in these two figures, it is clear that the simulations agree better with the experiments for the $\theta = 43^{\circ}$ angle. This is expected since jet noise peaks at shallower angles [87]. From Fig. 9b, we can also see that the effect of the wing is very small at $\theta = 43^{\circ}$. In contrast,



Figure 8: Velocity statistics in jet plume for the isolated and the installed jet. Isolated-1 (____), Installed-1 (____), Installed-2 (____), Experiments (O).

the far-field noise levels are considerably higher for the installed nozzle at the $\theta = 90^{\circ}$ angle, especially for St < 2 (Fig. 9a). These results are explained by the fact that jet installation noise has a dipole directivity that peaks in the $\theta = 90^{\circ}$ and 270° direction [88].

Figures 9a and 9b show that the simulation of the isolated nozzle tends to 782 over-predict the corresponding experimental results, especially for the higher 783 frequencies. We believe these discrepancies can be partly explained by the 784 under-resolved boundary layer inside the nozzle. The resolution (mesh and 785 polynomial degree) might also be too low for the isolated nozzle. In fact, if 786 we compare the spectra obtained with the P = 2 and P = 3 simulation of 787 the installed nozzle, we see that the simulation with the higher polynomial 788 degree (Installed-2) gives lower noise levels for higher frequencies. Since jet 789 installation noise mostly contributes to the lower frequencies, it is likely that 790 the noise levels for the isolated jet would come down further if P = 3 was 791 used. Unfortunately, this turned out to be outside the scope of this paper. 792

Despite the improvement seen when going from P = 2 to P = 3 for 793 the installed nozzle, there are still some notable discrepancies in the far-794 field spectra for the highest and lowest frequencies. The discrepancies seen 795 for the highest frequencies are likely caused by the same error sources as 796 for the isolated nozzle, e.g., insufficient modeling of the boundary layer and 797 resolution in the jet plume. Also, since the microphones are located below 798 the wing, the errors at the highest frequencies are amplified due to reflections 799 against the pressure side of the wing. With regards to the lowest frequencies, 800 on the other hand, the effect of the resolution along the wing span and size 801 of the Ffowcs Williams - Hawkings integration surface could be potential 802



Figure 9: Power Spectral Density in the far-field. Isolated-1 (____), Installed-1 (____), Installed-2 (____), Experiments for isolated jet (____), Experiments for installed jet (____).

sources of error. The effect of these should be investigated in future work.

804 4.2. Scaling and performance

The scalability of Nektar++ has been demonstrated on a number of parallel computers. In this case, strong scaling results of the compressible flow solver in Nektar++ on ARCHER2 and JUWELS supercomputers are presented. The scaling plots are generated based on the time-integration computational time (whole application except initialization and I/O). This approach is taken since the majority of time is spent on time-integration in a production run.

The DG method and 2nd order implicit time integration scheme introduced in sections 2.2.1 and 2.2.2, respectively, are considered for the scaling. The installed round nozzle case presented in sections 4.1.1 to 4.1.3 is used. This test has been run for 100 time steps on a mesh of approximately 5 mil⁸¹⁶ lion elements. Two polynomial orders, P = 2 and P = 3, have been tested. ⁸¹⁷ In total, this case contains approximately 54 million DOF for P = 2 and 109 ⁸¹⁸ million DOF for P = 3.

The scaling tests on ARCHER2 and JUWELS are presented in Fig. 10. 819 Tables including the timings of interest on ARCHER2 and JUWELS are 820 also presented in Table 3 and Table 4, respectively. It can be seen from 821 these figures that the scalings are done between 2,560 and 140,800 cores on 822 ARCHER2 and between 2,400 and 28,800 cores on JUWELS. ARCHER2 has 823 128 cores (2x AMD Zen2 Rome EPYC 7742 CPUs and 64 cores per CPU) 824 and 256 GB per node, while JUWELS has 48 cores (2x Intel Xeon Platinum 825 8168 Skylake CPUs and 24 cores per CPU) and 96 GB per node. We be-826 lieve that the higher memory per node on ARCHER2 allowed us to use a 827 higher number of nodes compared to JUWELS. In Fig. 10 we observe a pro-828 nounced super-linear scaling on ARCHER2 probably due to a larger number 829 of cache misses for small job sizes. This probably means that Nektar++ 830 is not cache optimized [89]. We can conclude from the scaling results that 831 the optimal number of DOF per core is between 650 and 850. Below that, 832 parallel efficiency decreases. 833

834 5. Conclusions

In this paper, we have presented a computational framework for accurately predicting jet noise. Our framework utilizes the spectral hp approach of Nektar++ to simulate turbulent jets. To compute the noise propagation from the jet vicinity to the far-field, we employed the Ffowcs Williams -Hawkings method, which is implemented in the Antares library. The primary

Num. nodes	Num. cores	CPU time $P=2$	CPU time P=3
20	2560	1309.62	3083.04
50	6400	410.658	1066.50
100	12800	191.250	496.383
150	19200	123.885	308.756
200	25600	91.6463	221.697
250	32000	78.5870	176.408
300	38400	60.2471	140.109
350	44800	52.2270	120.575
400	51200	44.9938	105.134
450	57600	40.8790	95.1388
500	64000	37.6927	84.2975
550	70400	32.3210	-
600	76800	31.6789	70.1267
650	83200	29.4066	-
700	89600	30.5700	61.5184
800	102400	31.6817	52.5133
850	108800	-	49.4919
900	115200	-	47.9693
1000	128000	-	43.7558
1100	140800	-	53.4228

Table 3: Time-integration CPU time of the scaling runs done on ARCHER2.



Figure 10: Strong scaling for the installed round nozzle case on ARCHER2, between 2,560 and 140,800 cores, and on JUWELS, between 2,400 and 28,800 cores, for polynomial orders P = 2 and P = 3.

objective of our work is to provide a comprehensive description of the computational framework developed, highlighting the integration of Nektar++
with Antares to enable turbulent jet simulations and reliable predictions of
jet noise.

To enable high-order simulations for turbulent jets and noise predictions, several important methodological aspects were discussed. Firstly, the generation of high-order meshes that accurately conform to the underlying CAD geometry was described. Additionally, the importance of establishing compatible initial and boundary conditions for high-order simulations was highlighted. In terms of post-processing, we presented a data acquisition strategy focused on probing the solution. As part of this strategy, we extended the

Num. nodes	Num. cores	CPU time $P=2$	CPU time $P=3$
50	2400	919.817	1811.14
100	4800	435.883	870.604
200	9600	192.249	416.621
500	24000	64.8598	166.230
600	28800	65.9233	127.540

Table 4: Time-integration CPU time of the scaling runs done on JUWELS.

history point filter in Nektar++ to store points in HDF5 format. Finally, we
addressed the process of generating the time history required for computing
noise propagation using Antares.

To showcase the computational framework, two configurations of a single 854 stream subsonic jet were computed. The first configuration focuses on an 855 isolated jet nozzle, while the second configuration incorporates a NACA4415 856 airfoil positioned near the nozzle exit. The numerical simulations were com-857 pared with experimental data provided by the University of Southampton. 858 The numerical aerodynamic results demonstrated very good agreement with 859 the experimental data. Small discrepancies were observed in the centreline 860 of the jet and further downstream. These discrepancies may be explained 861 by the under-resolved boundary layer and the inclusion of only a short part 862 of the conical nozzle. In terms of far-field results, some discrepancies were 863 observed between the simulations and the experiments. In particular, the 864 simulations tend to over-predict the corresponding experimental results for 865 the isolated nozzle, mainly for the highest frequencies. For the installed noz-866 zle, discrepancies in the far-field spectra were observed for the highest and 867

lowest frequencies. For the highest frequencies, the discrepancies were attributed to the insufficient modeling of the boundary layer and resolution in the jet plume in both cases. On the other hand, the discrepancies in the lowest frequencies for the installed nozzle could potentially be caused by a relatively low resolution along the wing span and size of the Ffowcs Williams - Hawkings integration surface around the wing. These source errors will be investigated in future work.

Recent developments in computational performance have resulted in sig-875 nificant speed up for highly computationally intensive simulations. One note-876 worthy development is the use of SIMD vectorization to calculate the diffu-877 sion operator and preconditioner within the compressible implicit solver in 878 Nektar++. Furthermore, the scalability of the Nektar++ framework has 879 been demonstrated on ARCHER2 and JUWELS supercomputers. The scal-880 ing results indicate linear scaling up to more than 100,000 cores for the 881 ARCHER2 supercomputer. 882

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