



A Performance and Energy Study of the Hyperbolic PDE Solver Engine ExaHyPE

Master's Thesis in Computational Science and Engineering

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September 6, 2016

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I hereby declare that this thesis is entirely the result of my own work except where otherwise indicated. I have only used the resources given in the list of references.

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Abstract

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Chapter 1

Introduction

- Challenges of exascale
 - Power:
 - * cite RAPL Memory: Datacenters are already overprovisioned (worst-case average instead of peak power)
 - * cite Energy-aware PAM for HPC: Methods: Direct measurement using specific hardware components (expensive, non-standardized, complex, per package). Software estimates or modeling (inaccurate, finer insight, cite rapl memory: For RAM: $\sigma = 1.1\text{percent}$), hybrid
- The ExaHyPE project (numerics, resilience, profiling) as an answer
- On the importance of profiling and performance measuring

Theoretical Context

2.1 A D-dimensional ADER-DG Scheme With MUSCL-Hancock a Posteriori Subcell Limiting for Non-linear Hyperbolic Balance Laws

2.1.1 Introduction

As stated already in the introduction, the ExaHyPE project is concerned with building an engine for simulating problems that can be formulated in terms of a hyperbolic balance law (HBL). Leaving all of the high performance computing (HPC) components aside, the very heart of the engine therefore comprises of an embedded numerical scheme for solving partial differential equations (PDEs) of HBL type.

Solving PDEs of this kind is of great interest in practice and has been a topic of active research for about a century¹. A comprehensive overview on schemes that have been proposed over the years can be found in [2]. Two particularly challenging aspects of such simulations are

1. the accurate simulation of HBL problems over long periods of time, especially when facing (prescribed or spontaneously arising) discontinuities (“shocks”) or stiff source terms leading to stability issues in space and time, respectively, and,
2. the inherent data access patterns the numerical schemes inherit from the PDE that make it challenging to avoid excessive communication

¹The British scientist and Durham University alumnus Lewis Fry Richardson is considered to be one of the founding fathers of computational fluid dynamics (CFD). In 1922 he published a book in which he presented a method for weather forecasting based on the solution of differential equations, in a time when most computations were still done by human “computers”. See [1] for a second edition of the book published in 2007 which puts the work into a contemporary context and emphasizes how modern weather forecasting is still based on Richardson’s ideas.

and to achieve high arithmetic density, two key drivers for good performance on modern distributed HPC systems.

In ExaHyPE a state of the art method called Arbitrary High Order Derivatives Discontinuous Galerkin (ADER-DG) is employed together with a posteriori subcell limiting based on the robust, second-order MUSCL-Hancock finite volume method (FVM) scheme (see [3] for details). As the name implies, the Discontinuous Galerkin framework with high-order local polynomials ansatz functions acts as the theoretical foundation of the method (see [4, 5] for more details on the underlying theory). In the following sections we will step by step derive the complete scheme and emphasize how it addresses the challenges stated above. The first part on unlimited ADER-DG is based to varying degree on work presented in [6] with three important additional contributions:

1. The scheme is presented in a more general form for systems involving V quantities in D spatial dimensions.
2. We employ index notation and simplify the equations up to a point where the resulting mathematical formulae can easily be mapped to a computer programming language, in particular it is in direct agreement with FORTRAN code used by Dumbser et al. to generate the numerical results presented in [7].
3. We extend the formulation to include a posteriori subcell limiting, introduce projection and reconstruction operations to equidistant subgrids and review the MUSCL-Hancock FVM scheme.

The chapter is set up as follows: First, we will begin with some remarks on the notation employed and state the problem at hand in its general form. Second, we will derive an element local weak formulation and approximate it with respect to finite-dimensional function spaces. After introducing reference coordinates, corresponding mappings and orthogonal bases for the function spaces involved, we employ a predictor-corrector approach to arrive at a fully-discrete method that is of arbitrarily high-order in both space and time.

2.1.2 Notation

Before we begin deriving the numerical schemes let us quickly introduce the following set of rules on how to depict common mathematical objects and operations:

- Vectors and vector-valued functions, i.e. first-order tensors, will be denoted by bold, lower case letters, e.g. \mathbf{x} , $\mathbf{u}(\mathbf{x}, t)$.
- Matrices or matrix-valued functions, i.e. second-order tensors, will be denoted by bold, upper case letters, e.g. \mathbf{K} , $\mathbf{F}(\mathbf{x}, t)$.

2.1. A D-dimensional ADER-DG Scheme With MUSCL-Hancock a Posteriori Subcell Limiting for Non-linear Hyperbolic Balance Laws

- Higher order tensors are always denoted as bold, lower-case letters with a “hat” on top, e.g. $\hat{\mathbf{u}}^{K,i}$, $\hat{\mathbf{q}}^{K,i}$. Note, however, that the opposite is not true².
- To avoid confusion in case we deal with tensors for which a superscript or a subscript is “part of its name” or where this indicates membership in a set, similar to the convention in many programming languages we denote “accesses” into the tensor by square brackets. An example illustrating the advantage of this notation based on common naming conventions in literature could be the following: Let $\{\hat{\mathbf{u}}_h^{K,i}\}_{K \in \mathcal{K}_h, i \in \{0,1,\dots,I-1\}}$ be the set of vector-valued functions from an appropriate Hilbert space which are defined locally on a cell $K \in \mathcal{K}_h$ and in a time interval $[t_i, t_i + \Delta t_i]$, $i \in \{0,1,\dots,I\}$. If we now want to access the v -th component of $\hat{\mathbf{u}}^{K,i}$, we write

$$\left[\hat{\mathbf{u}}_h^{K,i}\right]_v \quad (2.1)$$

instead of

$$\hat{\mathbf{u}}_{h,v}^{K,i} \quad (2.2)$$

In this way it is absolutely clear that K , i and h are “part of the name” and that v is an index used to access an element in the tensor. However in case an expression is absolutely unambiguous, for the sake of brevity we will often omit the square brackets. Most prominently we write

$$\frac{\partial}{\partial x_d} \quad (2.3)$$

instead of

$$\frac{\partial}{\partial [x]_d}. \quad (2.4)$$

- Throughout the thesis we use index notation following the Einstein summation convention whenever possible. This means that if an index within a product expression is repeated exactly once, this implies summation over the whole value range of this index. The standard inner product of two vectors $\mathbf{x}, \mathbf{y} \in \mathbb{R}^N$ can then be written as

$$\langle \mathbf{x}, \mathbf{y} \rangle = \sum_{n=0}^{N-1} [\mathbf{x}]_n [\mathbf{y}]_n = [\mathbf{x}]_n [\mathbf{y}]_n (= [\mathbf{x}]_m [\mathbf{y}]_m) \quad (2.5)$$

²In general the dimensionality of an object will always be obvious from its indices. Since we will just allow scalar and vector-valued indices only this distinction is of critical importance.

and for unambiguous cases like above

$$\langle \mathbf{x}, \mathbf{y} \rangle = \sum_{n=0}^{N-1} \mathbf{x}_n \mathbf{y}_n := \mathbf{x}_n \mathbf{y}_n (= \mathbf{x}_m \mathbf{y}_m). \quad (2.6)$$

Such indices are called dummy indices in the sense that as illustrated in the example above it does not matter if the index is named n or m . Sometimes free indices, i.e. indices that are not dummy indices, appear twice in a term as a result of some algebraic manipulation. In these cases we will explicitly state that summation over the index is not intended, unless it is obvious e.g. from the left-hand side of the equation that the index can only be a free index. We furthermore always give an explicit range for free indices. See [8] for more details on index notation, its advantages and disadvantages as well as a more formal definition.

In addition to increased brevity, index notation allows for less ambiguities compared to classical vector notation, simplifies derivation of identities from tensor calculus and if done carefully the resulting formulae can be mapped conveniently to loops in low-level programming languages such as C or FORTRAN.

- To keep all derivations dimension-agnostic, we define accesses into tensors using vector indices as follows: Let $\hat{\mathbf{u}} \in \mathbb{R}^{I_1 \times I_2 \times \dots \times I_D}$ be a tensor of order $D \in \mathbb{N}$ with $I_d \in \mathbb{N}_0$ for all $d \in \{0, 1, \dots, D-1\} := \mathcal{D}$. Let furthermore $i_d \in \{0, 1, \dots, I_d-1\}$ for all $d \in \mathcal{D}$ and $\mathbf{i} \in \mathbb{N}_0^D$, $[\mathbf{i}]_d = i_d$ for $d \in \mathcal{D}$ a vector of indices. Then we define

$$[\hat{\mathbf{u}}]_{\mathbf{i}} = [\hat{\mathbf{u}}]_{[i_0, i_1, \dots, i_{D-1}]} = [\hat{\mathbf{u}}]_{i_0, i_1, \dots, i_{D-1}}. \quad (2.7)$$

If we only provide a vector of $D-1$ indices, i.e.

$$[\hat{\mathbf{u}}]_{[i_0, i_1, \dots, i_{D-2}]} \quad (2.8)$$

we obtain a vector of length I_{D-1} . If we only provide $D-2$ indices we obtain a matrix with I_{D-2} rows and I_{D-1} columns. In general if we provide $d \in \{0, 1, \dots, D\}$ indices we obtain a tensor of order $D-d$.

- In the style of numerical computing environments such as MATLAB or Octave [9] we define the following shorthand notation for sequences of consecutive integral numbers:

$$j:k := \begin{cases} \{j, j+1, \dots, k\} & \text{if } j \leq k \\ \{\} & \text{otherwise.} \end{cases} \quad (2.9)$$

- We can now define access into a vector $\mathbf{x} \in \mathbb{R}^N$ of length N via sequences as

$$[\mathbf{x}]_{j:k} := \left[[\mathbf{x}]_j, [\mathbf{x}]_{j+1}, \dots, [\mathbf{x}]_k \right] \quad (2.10)$$

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for $j \leq k$ and $j, k \in 0:N-1$, which for unambiguous cases as above is equal to the definition

$$\mathbf{x}_{j:k} := [x_j, x_{j+1}, \dots, x_k]. \quad (2.11)$$

Together with implicit set and vector concatenation we can then write

$$[\mathbf{x}]_{\{0:k-1, k+1:N-1\}} = \mathbf{x}_{\{0:k-1, k+1:N-1\}} \quad (2.12)$$

for $k \in 0:N-1$ to denote the vector of length $N-1$ that contains all values of the original vector \mathbf{x} but the k -th component. Furthermore

$$[[\mathbf{x}]_{0:k-1}, x', [\mathbf{x}]_{k+1:N-1}] = [\mathbf{x}_{0:k-1}, x', \mathbf{x}_{k+1:N-1}] \quad (2.13)$$

denotes the vector of length N whose components are equal to the ones of \mathbf{x} apart from the k -th one, which we have replaced by the scalar $x' \in \mathbb{R}$.

2.1.3 Hyperbolic Balance Laws

A D -dimensional balance law in a system with V quantities is described mathematically by a partial differential equation (PDE) of the form

$$\frac{\partial}{\partial t} [\mathbf{u}(\mathbf{x}, t)]_v + \frac{\partial}{\partial x_d} [\mathbf{F}(\mathbf{u}(\mathbf{x}, t))]_{vd} = [\mathbf{s}(\mathbf{u}(\mathbf{x}, t))]_v \text{ on } \Omega \times [0, T] \quad (2.14)$$

together with initial conditions

$$[\mathbf{u}(\mathbf{x}, 0)]_v = [\mathbf{u}_0(\mathbf{x})]_v \quad \forall \mathbf{x} \in \Omega, \quad (2.15)$$

and boundary conditions

$$[\mathbf{u}(\mathbf{x}, t)]_v = [\mathbf{u}_B(\mathbf{x}, t)]_v \quad \forall \mathbf{x} \in \partial\Omega, t \in [0, T], \quad (2.16)$$

for all $v \in \mathcal{V}$, where we define the index set $\mathcal{V} := \{1, 2, \dots, V\}$. $[0, T]$ is the time interval of interest and $\Omega \subset \mathbb{R}^D$ denotes the spatial domain. The function $\mathbf{F} : \mathbb{R}^V \rightarrow \mathbb{R}^{V \times D}$, $\mathbf{u} \mapsto \mathbf{F}(\mathbf{u}) = [f_1(\mathbf{u}), f_2(\mathbf{u}), \dots, f_D(\mathbf{u})]$ is called the flux function. For the problem to be hyperbolic we require that all Jacobian matrices $\mathbf{A}_d(\mathbf{u})$, $d \in \{0, 1, \dots, D-1\} := \mathcal{D}$, defined as

$$[\mathbf{A}_d]_{ij} = \frac{\partial [f_d]_i}{\partial u_j}, \quad (2.17)$$

have D real eigenvalues in each admissible state $\mathbf{u} \in \mathbb{R}^V$.

2.1.4 Space and Time Discretization

Let \mathcal{K}_h be a quadrilateral partition of Ω , i.e.

$$K \cap J = \emptyset \forall K, J \in \mathcal{K}_h, K \neq J, \quad (2.18)$$

$$\bigcup_{K \in \mathcal{K}_h} K = \Omega. \quad (2.19)$$

For the index set $\mathcal{I} := \{0, 1, \dots, I-1\}$ let $\{t_i\}_{i \in \mathcal{I}}$ be an I -fold partition of the time interval $[0, T]$ such that

$$0 = t_0 < t_1 < \dots < t_I = T. \quad (2.20)$$

For $i \in \mathcal{I}$ we furthermore define

$$\Delta t_i = t_{i+1} - t_i, \quad (2.21)$$

so that the subinterval $[t_i, t_{i+1}]$ can be written as $[t_i, t_i + \Delta t_i]$.

Without loss of generality we can solve the original PDE (2.14) on $\Omega \times [0, T]$ simply by solving the PDE locally for each element $K \in \mathcal{K}_h$ in the time interval $[t_0, t_0 + \Delta t_0]$ and then proceeding to the next time interval until we have reached the final time T . This gives rise to an element-local formulation on a subinterval in time which we will focus on in the following.

2.1.5 Element-local Weak Formulation

Let $L^2(\Omega)^V$ be the space of vector-valued, square-integrable functions on Ω , i.e.

$$L^2(\Omega)^V = \left\{ \mathbf{w} : \Omega \rightarrow \mathbb{R}^V \mid \int_{\Omega} \|\mathbf{w}\|^2 dx < \infty \right\}. \quad (2.22)$$

Let $\mathbf{w} \in L^2(\Omega)^V$ be a spatial test function. Multiplication of the original PDE (2.14) and integration over a space-time cell $K \times [t_i, t_i + \Delta t_i]$ yields an element-local weak formulation of the problem,

$$\begin{aligned} & \int_{t_i}^{t_i + \Delta t_i} \int_K \frac{\partial}{\partial t} [\mathbf{u}]_v [\mathbf{w}]_v dx dt + \int_{t_i}^{t_i + \Delta t_i} \int_K \frac{\partial}{\partial x_d} [\mathbf{F}(\mathbf{u})]_{vd} [\mathbf{w}]_v dx dt = \\ & \int_{t_i}^{t_i + \Delta t_i} \int_K [\mathbf{s}(\mathbf{u})]_v [\mathbf{w}]_v dx dt, \end{aligned} \quad (2.23)$$

which we require to hold for all $v \in \mathcal{V}$, $\mathbf{w} \in L^2(\Omega)^V$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

Integration by parts of the spatial integral in the second term yields

$$\begin{aligned} & \int_K \frac{\partial}{\partial x_d} [\mathbf{F}(\mathbf{u})]_{vd} [\mathbf{w}]_v dx = \\ & \int_K \frac{\partial}{\partial x_d} \left([\mathbf{F}(\mathbf{u})]_{vd} [\mathbf{w}]_v \right) dx - \int_K [\mathbf{F}(\mathbf{u})]_{vd} \frac{\partial}{\partial x_d} [\mathbf{w}]_v dx. \end{aligned} \quad (2.24)$$

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Application of the divergence theorem to the first term on the right-hand side of (2.24) yields

$$\int_K \frac{\partial}{\partial x_d} \left([F(\mathbf{u})]_{vd} [w]_v \right) d\mathbf{x} = \int_{\partial K} [F(\mathbf{u})]_{vd} [w]_v [\mathbf{n}]_d ds(\mathbf{x}), \quad (2.25)$$

where $\mathbf{n} \in \mathbb{R}^D$ is the unit-length, outward-pointing normal vector at a point \mathbf{x} on the surface of K which we denote by ∂K .

Inserting eqs. (2.24) and (2.25) into eq. (2.23) yields the following more favorable element-local weak formulation of the original equation (2.14):

$$\begin{aligned} & \int_{t_i}^{t_i+\Delta t_i} \int_K \frac{\partial}{\partial t} [\mathbf{u}]_v [w]_v d\mathbf{x} dt - \int_{t_i}^{t_i+\Delta t_i} \int_K [F(\mathbf{u})]_{vd} \frac{\partial}{\partial x_d} [w]_v d\mathbf{x} dt + \\ & \int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} [F(\mathbf{u})]_{vd} [w]_v [\mathbf{n}]_d ds(\mathbf{x}) dt = \int_{t_i}^{t_i+\Delta t_i} \int_K [s(\mathbf{u})]_v [w]_v d\mathbf{x} dt. \end{aligned} \quad (2.26)$$

Again we require the weak formulation to hold for all $v \in \mathcal{V}$, $w \in L^2(\Omega)^V$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

2.1.6 Restriction to Finite-Dimensional Function Spaces

To discretize eq. (2.26) we need to impose the restriction that both test and ansatz functions come from a finite-dimensional function space. First, let $\mathcal{Q}_N(K)^V$ and $\mathcal{Q}_N(K \times [t_i, t_i + \Delta t_i])^V$ be the space of vector-valued, multivariate polynomials of degree less or equal than N in each variable on K and $K \times [t_i, t_i + \Delta t_i]$, respectively. We can then define the following finite-dimensional function spaces:

- For spatial functions we define

$$\mathbb{W}_h = \left\{ w_h \in L^2(\Omega)^V \mid w_h|_K := w_h^K \in \mathcal{Q}_N(K)^V \forall K \in \mathcal{K}_h \right\}. \quad (2.27)$$

- For space-time functions on the time subinterval $[t_i, t_i + \Delta t_i]$, $i \in \mathcal{I}$ we define

$$\begin{aligned} \tilde{\mathbb{W}}_h^i = & \left\{ \tilde{w}_h^i \in L^2(\Omega \times [t_i, t_i + \Delta t_i]) \mid \right. \\ & \left. \tilde{w}_h^i|_K := \tilde{w}_h^{K,i} \in \mathcal{Q}_N(K \times [t_i, t_i + \Delta t_i]) \forall K \in \mathcal{K}_h \right\}. \end{aligned} \quad (2.28)$$

2. THEORETICAL CONTEXT

Replacing w by $w_h \in \mathbb{W}_h$ and u by $\tilde{u}_h^i \in \tilde{\mathbb{W}}_h^i$ in eq. (2.26), i.e. restricting ourselves to test and ansatz functions from finite-dimensional function spaces, yields an approximation of the weak formulation,

$$\begin{aligned} & \int_{t_i}^{t_i+\Delta t_i} \int_K \frac{\partial}{\partial t} [\tilde{u}_h^{K,i}]_v [w_h^K]_v dxdt - \int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} [F(\tilde{u}_h^{K,i})]_{vd} \frac{\partial}{\partial x_d} [w_h^K]_v dxdt + \\ & \int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} [\mathcal{G}(\tilde{u}_h^{K,i}, \tilde{u}_h^{K',i}, n)]_v [w_h^K]_v ds(x)dt = \\ & \int_{t_i}^{t_i+\Delta t_i} \int_K [s(\tilde{u}_h^{K,i})]_v [w_h^K]_v dxdt, \end{aligned} \quad (2.29)$$

which now has to hold for all $w_h \in \mathbb{W}_h$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$. Since for a cell $K \in \mathcal{K}_h$ and one of its Voronoi neighbors $K' \in V(K)$ in general it holds that

$$\tilde{u}_h^{K,i}(x^*) \neq \tilde{u}_h^{K',i}(x^*) \quad (2.30)$$

for $x^* \in K \cap K'$, i.e. \tilde{u}_h^i is double-valued at the interface between K and K' , in order to compute the surface integral we need to introduce the numerical flux function $\mathcal{G}(\tilde{u}_h^{K,i}, \tilde{u}_h^{K',i}, n)$. The numerical flux at a position $x^* \in K \cap K'$ on the interface is obtained by (approximately) solving a Riemann problem in normal direction.

Excursus: The Riemann Problem

Let x^* be a point on the interface ∂K between a cell $K \in \mathcal{K}_h$ and its Voronoi neighbor $K' \in V(K)$ and let n be the outward pointing unit normal vector at this point. Then to obtain the numerical flux we need to solve the initial boundary value problem (“Riemann problem”)

$$\frac{\partial}{\partial t} [g]_v + \sum_{d=1}^D \frac{\partial}{\partial x_d} [F(g)]_{vd} [n]_d = 0 \quad (2.31)$$

along the line $x = x^* + \alpha n$ for $\alpha \in \mathbb{R}$ with discontinuous initial conditions

$$g(x^* + \alpha n, 0) = \begin{cases} \tilde{u}_h^{K,i}|_{x^*} & \text{if } \alpha < 0 \\ \tilde{u}_h^{K',i}|_{x^*} & \text{if } \alpha > 0. \end{cases} \quad (2.32)$$

We then evaluate the similarity solution $\tilde{g}(\alpha/t)$ of the problem and define

$$\left[\mathcal{G}(\tilde{u}_h^{K,i}, \tilde{u}_h^{K',i}, n) \right]_v := \left[\tilde{g}|_0 \right]_v. \quad (2.33)$$

For an extensive overview on state of the art approximate Riemann solvers see [2].

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Continuing with eq. (2.29), integration by parts in time of the first term and observing that w_h is constant in time yields the following one-step update scheme for the cell-local time-discrete solution $\tilde{u}_h^{K,i}$:

$$\begin{aligned} \int_K \left[\tilde{u}_h^{K,i} \Big|_{t_i+\Delta t_i} \right]_v \left[w_h^K \right]_v dx &= \int_K \left[\tilde{u}_h^{K,i} \Big|_{t_i} \right]_v \left[w_h^K \right]_v dx + \\ &\quad \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F(\tilde{u}_h^{K,i}) \right]_{vd} \frac{\partial}{\partial x_d} \left[w_h^K \right]_v dx dt - \\ &\quad \int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} \left[\mathcal{G}(\tilde{u}_h^{K,i}, \tilde{u}_h^{K+i}, n) \right]_v \left[w_h^K \right]_v ds(x) dt + \\ &\quad \int_{t_i}^{t_i+\Delta t_i} \int_K \left[s(\tilde{u}_h^{K,i}) \right]_v \left[w_h^K \right]_v dx dt. \end{aligned} \quad (2.34)$$

Again we require eq. (2.34) to hold for all $v \in \mathcal{V}$, $w_h \in \mathbb{W}_h$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$. Note, however, that the scheme is incomplete, since we only know $\tilde{u}_h^i|_t$ at the discrete time steps $t \in \{t_i, t_i + \Delta t_i\}$, not within the open interval, i.e. for $t \in (t_i, t_i + \Delta t_i)$. As commonly done in a DG framework we therefore proceed by replacing \tilde{u}_h on the interval $(t_i, t_i + \Delta t_i)$ by an approximation $\tilde{q}_h^i \in \tilde{\mathbb{W}}_h^i$ which we call space-time predictor.

2.1.7 Space-time Predictor

To derive a procedure to compute the space-time predictor $\tilde{q}_h^i \in \tilde{\mathbb{W}}_h^i$ we again start from the original PDE (2.14), but this time we do not use a spatial test function $w_h \in \mathbb{W}_h$, but a space-time test function $\tilde{w}_h^i \in \tilde{\mathbb{W}}_h^i$. If we furthermore replace the solution u by the space-time predictor $\tilde{q}_h^i \in \tilde{\mathbb{W}}_h^i$, integrate over the space-time element $K \times [t_i, t_i + \Delta t_i]$ and apply the divergence theorem analogously to eq. (2.25) we obtain the following relation:

$$\begin{aligned} &\int_{t_i}^{t_i+\Delta t_i} \int_K \frac{\partial}{\partial t} \left[\tilde{q}_h^{K,i} \right]_v \left[\tilde{w}_h^{K,i} \right]_v dx dt - \\ &\quad \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F(\tilde{q}_h^{K,i}) \right]_{vd} \frac{\partial}{\partial x_d} \left[\tilde{w}_h^{K,i} \right]_v dx dt + \\ &\quad \int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} \left[\mathcal{G}(\tilde{q}_h^{K,i}, \tilde{q}_h^{K+i}, n) \right]_v \left[\tilde{w}_h^{K,i} \right]_v ds(x) dt = \\ &\quad \int_{t_i}^{t_i+\Delta t_i} \int_K \left[s(\tilde{q}_h^{K,i}) \right]_v \left[\tilde{w}_h^{K,i} \right]_v dx dt. \end{aligned} \quad (2.35)$$

We require eq. (2.35) to hold for all $v \in \mathcal{V}$, $\tilde{w}_h^i \in \tilde{\mathbb{W}}_h^i$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

The assumption that the solution is balanced, i.e. that there is no net inflow or outflow for cells $K \in \mathcal{K}_h$ allows us to drop the third term. Together with integration by parts in time applied to the first term this yields

$$\begin{aligned} & \int_K \left[\tilde{q}_h^{K,i} \Big|_{t_i+\Delta t_i} \right]_v \left[\tilde{w}_h^{K,i} \Big|_{t_i+\Delta t_i} \right]_v dx - \int_{t_i}^{t_i+\Delta t_i} \int_K \left[\tilde{q}_h^{K,i} \right]_v \frac{\partial}{\partial t} \left[\tilde{w}_h^{K,i} \right]_v dx dt = \\ & \int_K \left[\tilde{q}_h^{K,i} \Big|_{t_i} \right]_v \left[\tilde{w}_h^{K,i} \Big|_{t_i} \right]_v dx + \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F(\tilde{q}_h^{K,i}) \right]_{vd} \frac{\partial}{\partial x_d} \left[\tilde{w}_h^{K,i} \right]_v dx dt + \\ & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[s \left(\tilde{q}_h^{K,i} \right) \right]_v \left[\tilde{w}_h^{K,i} \right]_v dx dt, \end{aligned} \quad (2.36)$$

which we require to hold for all $v \in \mathcal{V}$, $\tilde{w}_h^i \in \tilde{W}_h^i$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$. In conjunction with the initial condition

$$\tilde{q}_h^{K,i} \Big|_{t_i} = \tilde{u}_h^{K,i} \quad (2.37)$$

and an initial guess

$$\tilde{q}_h^{K,i} \Big|_t = \tilde{u}_h^{K,i} \forall t \in (t_i, t_i + \Delta t_i] \quad (2.38)$$

this relation can be used as a fixed-point iteration to find the cell-local space-time predictor $\tilde{q}_h^{K,i}$.

In the following two sections we will introduce mappings from spatial elements K and space-time elements $K \times [t_i, t_i + \Delta t_i]$ to spatial and space-time reference cells as well as orthogonal bases for the spaces W_h and \tilde{W}_h^i . We will then insert these results into eq. (2.36) and derive a fully-discrete iterative method to compute the cell-local space-time predictor $\tilde{q}_h^{K,i}$.

2.1.8 Reference Elements and Mappings

Let $\hat{K} := [0, 1]^D$ be the spatial reference element and $\xi \in \hat{K}$ be a point therein. Let $[0, 1]$ be the reference time interval and $\tau \in [0, 1]$ be a point in reference time. We can then introduce the following mappings:

Spatial mappings: Let $K \in \mathcal{K}_h$ be a cell in global coordinates with extent Δx^K and “lower-left corner” P_K , more precisely that is

$$[\Delta x^K]_d = \max_{x \in K} [x]_d - \min_{x \in K} [x]_d \quad (2.39)$$

and

$$[P_K]_d = \min_{x \in K} [x]_d \quad (2.40)$$

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for $d \in \mathcal{D}$. We can then define a mapping

$$\mathcal{X}_K : \hat{K} \rightarrow K, \xi \mapsto \mathcal{X}_K(\xi) = x \quad (2.41)$$

via the relation

$$[x]_d = [\mathcal{X}_K(\xi)]_d = [P_K]_d + [\Delta x]_d [\xi]_d \quad (2.42)$$

for $d \in \mathcal{D}$ (i.e. no summation on d) and for all $x \in K$, $\xi \in \hat{K}$ and $K \in \mathcal{K}_h$.

Temporal mappings: Let $[t_i, t_i + \Delta t_i], i \in \mathcal{I}$ be an interval in global time. The mapping

$$\mathcal{T}_i : [0, 1] \rightarrow [t_i, t_i + \Delta t_i], \tau \mapsto \mathcal{T}_i(\tau) = t_i + \Delta t_i \tau = t \quad (2.43)$$

maps a point $\tau \in [0, 1]$ in reference time to a point $t \in [t_i, t_i + \Delta t_i]$ in global time for all $i \in \mathcal{I}$.

The inverse mappings, the Jacobian matrices and the Jacobi determinants of the mappings are given in the following:

Spatial mappings: The inverse spatial mappings

$$\mathcal{X}_K^{-1} : K \rightarrow \hat{K}, x \mapsto \mathcal{X}_K^{-1}(x) = \xi \quad (2.44)$$

are defined via the relation

$$[\xi]_d = [\mathcal{X}_K^{-1}(x)]_d = \frac{1}{[\Delta x^K]_d} ([x]_d - [P_K]_d) \quad (2.45)$$

for $d \in \mathcal{D}$ and for all $\xi \in \hat{K}$, $x \in K$ and $K \in \mathcal{K}_h$. The Jacobian of \mathcal{X}_K is found to be

$$\left[\frac{\partial \mathcal{X}_K}{\partial \xi} \right]_{dd'} = \frac{\partial [\mathcal{X}_K]_d}{\partial \xi_{d'}} = [\Delta x^K]_d \delta_{dd'}, \quad (2.46)$$

where $d, d' \in \mathcal{D}$ (i.e. no summation on d) and for all $K \in \mathcal{K}_h$. As usual $\delta_{dd'}$ denotes the Kronecker delta defined as

$$\delta_{dd'} = \begin{cases} 0 & \text{if } d \neq d' \\ 1 & \text{if } d = d'. \end{cases} \quad (2.47)$$

The Jacobi determinant of \mathcal{X}_K for $K \in \mathcal{K}_h$ then simply is

$$J_{\mathcal{X}_K} = \left\| \frac{\partial \mathcal{X}_K}{\partial \xi} \right\| = \prod_{d=1}^D [\Delta x^K]_d, \quad (2.48)$$

i.e. the determinant is constant with respect to all $\xi \in \hat{K}$.

Temporal mappings: The inverse temporal mappings are given as

$$\mathcal{T}_i^{-1} : [t_i, t_i + \Delta t_i] \rightarrow [0, 1], t \mapsto \mathcal{T}_i^{-1}(t) = \frac{t - t_i}{\Delta t_i} = \tau \quad (2.49)$$

for all $\tau \in [0, 1]$, $t \in [t_i, t_i + \Delta t_i]$ and $i \in \mathcal{I}$. In the trivial case of a one-dimensional mapping the Jacobian of \mathcal{T}_i is a scalar which in turn is its own determinant. One finds

$$\frac{d\mathcal{T}_i}{d\tau} = \Delta t_i = J_{\mathcal{T}_i} \quad (2.50)$$

which again is constant with respect to all $\tau \in [0, 1]$.

2.1.9 Orthogonal Bases for the Finite-dimensional Function Spaces

In section 2.1.6 we introduced finite-dimensional, cell-wise polynomial function spaces \mathbb{W}_h and $\tilde{\mathbb{W}}_h^i$ for spatial and space-time ansatz and test functions, respectively. On our way towards a fully discrete version of the relations (2.36) and (2.34) to obtain the space-time predictor and the solution at the next time step, respectively, we will now derive a set of functions that form bases for the two function spaces of interest. Following the approach presented by Dumbars et al. in [3], throughout the thesis we will use the set of Lagrange functions with nodes located at the roots of the Legendre polynomials and tensor products thereof. In the later chapters of this work it will become obvious why this particular choice is highly favorable. For the moment the two major reasons shall be stated as an outlook:

1. Numerical integration using the Gauss-Legendre method is simple and computationally cheap, since the function values at the Gauss-Legendre nodes are directly available as they are equal to the degrees of freedom representing the local polynomial.
2. The resulting bases are orthogonal, which in turn makes sure that the resulting DG-matrices exhibit a sparse block structure allowing computations to be carried out efficiently in a dimension-by-dimension manner.

Lagrange Interpolation

Let $f \in \mathbb{Q}_N([0, 1])$ be a polynomial of degree less or equal than N and for the index set $\mathcal{N} := \{0, 1, \dots, N\}$ let $\{\hat{\xi}_n\}_{n \in \mathcal{N}}$ be a set of distinct nodes in $[0, 1]$. Then the Lagrange interpolation of f ,

$$\hat{f}(\xi) = \sum_{n=0}^N L_n(\xi) f(\xi_n) \quad (2.51)$$

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with Lagrange functions

$$L_n(\xi) = \prod_{m=0, m \neq n}^N \frac{\xi - \hat{\xi}_m}{\hat{\xi}_n - \hat{\xi}_m} \quad (2.52)$$

is exact, i.e.

$$f(\xi) = \hat{f}(\xi) \quad \forall \xi \in [0, 1]. \quad (2.53)$$

Since therefore every polynomial $f \in \mathbb{Q}_N([0, 1])$ can be represented as a linear combination of the Legendre polynomials L_n , $n \in \mathcal{N}$, the set of functions $\{L_n\}_{n \in \mathcal{N}}$ is a basis of $\mathbb{Q}_N([0, 1])$.

The following observation is an important property of the Lagrange polynomials:

$$L_n(\hat{\xi}_{n'}) = \delta_{nn'}, \quad (2.54)$$

i.e. at each node $\hat{\xi}_n$ only L_n has value 1 and all other polynomials evaluate to 0.

Legendre Polynomials and Gauss-Legendre Integration

Let $P_0 : [-1, 1] \rightarrow \mathbb{R}, \xi \mapsto 1$ and $P_1 : [-1, 1] \rightarrow \mathbb{R}, \xi \mapsto \xi$ be the zeroth and the first Legendre polynomial, respectively. Then the $N + 1$ -th Legendre polynomial can be defined via the following recurrence relation:

$$P_{N+1}(\xi) = \frac{1}{N+1} ((2N+1)P_N(\xi) - nP_{N-1}(\xi)). \quad (2.55)$$

Let $\{\tilde{\xi}_n\}_{n \in \mathcal{N}}$ be the roots of the $N + 1$ -th Legendre polynomial L_{N+1} . Then $\{\hat{\xi}_n\}_{n \in \mathcal{N}}$ with

$$\hat{\xi}_n = \frac{1}{2}(\tilde{\xi}_n + 1) \quad (2.56)$$

are the roots of the $N + 1$ -th Legendre polynomial linearly mapped to the interval $[0, 1]$. In conjunction with a set of suitable weights $\{\hat{\omega}_n\}_{n \in \mathcal{N}}$ Gauss-Legendre integration can be used to integrate polynomials of degree up to $2N + 1$ over the interval $[0, 1]$ exactly, i.e.

$$\int_0^1 f(\xi) d\xi = \sum_{n=0}^N \hat{\omega}_n f(\hat{\xi}_n) \quad \forall f \in \mathbb{Q}_{2N+1}([0, 1]). \quad (2.57)$$

A Python notebook on how to find the nodes $\{\hat{\xi}_n\}_{n \in \mathcal{N}}$ and weights $\{\hat{\omega}_n\}_{n \in \mathcal{N}}$ can be found in appendix A.

Scalar-valued Basis Functions on the One-dimensional Reference Element

Let $\{\hat{\psi}_n\}_{n \in \mathcal{N}}$ be the set of $N + 1$ Lagrange polynomials with nodes at the roots of the $N + 1$ -th Legendre polynomial linearly mapped to the interval $[0, 1]$, i.e.

$$\hat{\psi}_n(x) = \sum_{n'=0}^N \frac{x - \hat{x}_{n'}}{\hat{x}_n - \hat{x}_{n'}} \quad (2.58)$$

for $n \in \mathcal{N}$. Since $\{\hat{\psi}_n\}_{n \in \mathcal{N}}$ are Lagrange polynomials and the roots $\{\hat{x}_n\}_{n \in \mathcal{N}}$ are distinct, the set is a basis of $\mathbb{Q}_N([0, 1])$. Since furthermore

$$\langle \hat{\psi}_n, \hat{\psi}_m \rangle_{L^2([0,1])} = \int_0^1 \hat{\psi}_n(x) \hat{\psi}_m(x) dx = \sum_{n'=0}^N \hat{w}'_n \hat{\psi}_n(\hat{x}_{n'}) \hat{\psi}_m(\hat{x}_{n'}) = \hat{w}_n \delta_{mn} \quad (2.59)$$

for all $m, n \in \mathcal{N}$ (i.e. no summation over n), the set is even an orthogonal basis of $\mathbb{Q}_N([0, 1])$ with respect to the L^2 -scalar product as defined above. In this derivation we used the fact that $\hat{\psi}_n \hat{\psi}_m$ has degree $2N$ and that Gauss-Legendre integration with $N + 1$ nodes is exact for polynomials up to degree $2N + 1$.

Scalar-valued Basis Functions on the Spatial Reference Element

For the vector-valued index set $\mathcal{N} := \{0, 1, \dots, N\}^D$ let us define the set of scalar-valued spatial basis functions $\{\hat{\phi}_n\}_{n \in \mathcal{N}}$ on $\hat{K} := [0, 1]^D$ as

$$\hat{\phi}_n(\xi) = \prod_{d=1}^D \hat{\psi}_{[n]_d}([\xi]_d) = \hat{\psi}_{[n]_d}([\xi]_d), \quad (2.60)$$

i.e. $\{\hat{\phi}_n\}_{n \in \mathcal{N}}$ is the tensor product of $\{\hat{\psi}_n\}_{n \in \mathcal{N}}$ and as such it is a basis of $\mathbb{Q}([0, 1]^D) = \mathbb{Q}(\hat{K})$. If we define

$$[\hat{\xi}_n]_d = \hat{\xi}_{[n]_d} \quad (2.61)$$

and

$$\prod_{d=1}^D \hat{\omega}_{[n]_d}, \quad (2.62)$$

for all $d \in \mathcal{V}$ and $n \in \mathcal{N}$, we furthermore observe that the basis is orthogonal with respect to the L^2 -scalar product, since

$$\begin{aligned} \langle \hat{\phi}_n, \hat{\phi}_m \rangle_{L^2(\hat{K})} &= \int_{\hat{K}} \hat{\phi}_n(\xi) \hat{\phi}_m(\xi) d\xi = \\ &= \sum_{n' \in \mathcal{N}} \left(\hat{\omega}_{n'} \hat{\phi}_n(\hat{\xi}_{n'}) \hat{\phi}_m(\hat{\xi}_{n'}) \right) = \hat{\omega}_n \delta_{nm} \end{aligned} \quad (2.63)$$

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for all $n, m \in \mathcal{N}$. The natural extensions of the Kronecker delta for vector-valued indices is defined as follows:

$$\delta_{nm} = \prod_{d=1}^D \delta_{[n]_d[m]_d} = \delta_{[n]_d[m]_d}. \quad (2.64)$$

Scalar-valued Basis Functions on the Space-time Reference Element

Analogous to the procedure illustrated above for the spatial reference element \hat{K} we can define a basis $\{\hat{\theta}_{nl}\}_{n \in \mathcal{N}, l \in \mathcal{N}}$ of $\mathbb{Q}_N(\hat{K} \times [0, 1])$ on the reference space-time element $\hat{K} \times [0, 1]$ as

$$\hat{\theta}_{nl}(\xi, \tau) = \hat{\phi}_n(\xi) \hat{\psi}_l(\tau), \quad (2.65)$$

which again is orthogonal, since

$$\left\langle \hat{\theta}_{nl}, \hat{\theta}_{mk} \right\rangle_{L^2(\hat{K} \times [0, 1])} = \int_0^1 \int_{\hat{K}} \hat{\theta}_{nl} \hat{\theta}_{mk} d\xi d\tau = \hat{\omega}_n \hat{\omega}_l \delta_{nm} \delta_{lk} \quad (2.66)$$

for all $n, m \in \mathcal{N}$ and $l, k \in \mathcal{N}$.

Vector-valued Basis Functions on the Spatial Reference Element

If we define $\{\hat{\phi}_{nv}\}_{n \in \mathcal{N}, v \in \mathcal{V}}$ as

$$\hat{\phi}_{nv} = \hat{\phi}_n e_v, \quad (2.67)$$

where e_v is the v -th unit vector, i.e.

$$[e_v]_{v'} = \delta_{vv'} \quad (2.68)$$

for $v, v' \in \mathcal{V}$. Since

$$\begin{aligned} \left\langle \hat{\phi}_{nv}, \hat{\phi}_{n'v'} \right\rangle_{L^2(\hat{K})^V} &= \int_{\hat{K}} [\hat{\phi}_{nv}]_j [\hat{\phi}_{n'v'}]_j d\xi = \\ &= ([e_v]_j [e_{v'}]_j) \int_0^1 \int_{\hat{K}} \hat{\phi}_n \hat{\phi}_{n'} d\xi = \hat{\omega}_n \delta_{nn'} \delta_{vv'} \end{aligned} \quad (2.69)$$

for all $n, n' \in \mathcal{N}$ and $v, v' \in \{1, 2, \dots, V\}$ the set is an orthogonal basis for $\mathbb{Q}_N(\hat{K})^V$.

Vector-valued Basis Functions on the Space-time Reference Element

The set $\{\hat{\theta}_{nlv}\}_{n \in \mathcal{N}, l \in \mathcal{N}, v \in \mathcal{V}}$ defined as

$$\hat{\theta}_{nlv}(\xi, \tau) = \hat{\theta}_{nl}(\xi, \tau) e_v = \hat{\phi}_n(\xi) \hat{\psi}_l(\tau) e_v \quad (2.70)$$

is a basis of $\mathcal{Q}_N(\hat{K} \times [0, 1])^V$. Since furthermore

$$\left\langle \hat{\theta}_{nlv}, \hat{\theta}_{n'l'v'} \right\rangle_{L^2(\hat{K} \times [0, 1])^V} = \int_0^1 \int_{\hat{K}} \left[\hat{\theta}_{nlv} \right]_j \left[\hat{\theta}_{n'l'v'} \right]_j d\hat{\xi} d\tau = \hat{\omega}_n \hat{\omega}_l \delta_{nn'} \delta_{ll'} \delta_{vv'}, \quad (2.71)$$

for all $n, n' \in \mathcal{N}$, $l, l' \in \mathcal{N}$ and $v, v' \in \mathcal{V}$, the set is an orthogonal basis with respect to the respective L^2 -scalar product.

2.1.10 Basis Functions in Global Coordinates

We can use the mappings derived in chapter 2.1.8 to map the basis functions to global coordinates. For the vector-valued basis functions on a spatial element K we obtain

$$\phi_{nv}^K(x) = \begin{cases} \left(\hat{\phi}_{nv} \circ \mathcal{X}_K^{-1} \right)(x) & \text{if } x \in K \\ 0 & \text{otherwise,} \end{cases} \quad (2.72)$$

and for the vector-valued basis functions on a space-time element $K \times [t_i, t_i + \Delta t_i]$ we have

$$\theta_{nlv}^{Ki}(x, t) = \begin{cases} \left(\hat{\theta}_{nlv} \circ \left(\mathcal{X}_K^{-1}, \mathcal{T}_i^{-1} \right) \right)(x, t) & \text{if } x \in K \text{ and } t \in [t_i, t_i + \Delta t_i] \\ 0 & \text{otherwise} \end{cases} \quad (2.73)$$

for $n \in \mathcal{N}$, $l \in \{0, 1, \dots, N\}$ as well as $v \in \mathcal{V}$ and for all $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

2.1.11 A Fully-discrete Iterative Method for the Space-time Predictor

We recall relation (2.38) for the space-time predictor. Plugging in the initial condition (2.37) yields

$$\begin{aligned} & \int_K \left[\tilde{q}_h^{K,i} \Big|_{t_i+\Delta t_i} \right]_j \left[\tilde{w}_h^{K,i} \Big|_{t_i+\Delta t_i} \right]_j dx - \int_{t_i}^{t_i+\Delta t_i} \int_K \left[\tilde{q}_h^{K,i} \right]_j \frac{\partial}{\partial t} \left[\tilde{w}_h^{K,i} \right]_j dx dt = \\ & \int_K \left[\tilde{u}_h^{K,i} \Big|_{t_i} \right]_j \left[\tilde{w}_h^{K,i} \Big|_{t_i} \right]_j dx + \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F(\tilde{q}_h^{K,i}) \right]_{jk} \frac{\partial}{\partial x_k} \left[\tilde{w}_h^{K,i} \right]_j dx dt + \\ & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[s(\tilde{q}_h^{K,i}) \right]_j \left[\tilde{w}_h^{K,i} \right]_j dx dt, \end{aligned} \quad (2.74)$$

which we require to hold for all $\tilde{w}_h \in \tilde{W}_h$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

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Making use of the bases we derived in the previous section the cell-local space-time predictor $\hat{q}_h^{K,i}$ can be represented by a tensor of coefficients $\hat{q}^{K,i}$ (“degrees of freedom”) as follows:

$$\hat{q}_h^{K,i} = \left[\hat{q}^{K,i} \right]_{nlv} \theta_{nlv}^{Ki}. \quad (2.75)$$

The initial condition $\hat{u}_h^{K,i} \Big|_{t_i}$ can be represented as

$$\hat{u}_h^{K,i} \Big|_{t_i} = \left[\hat{u}^{K,i} \right]_{nv} \phi_{nv}^K, \quad (2.76)$$

where

$$\left[\hat{u}^{K,i} \right]_{nv} = \left[\hat{u}_h^{K,i} \Big|_{(\mathcal{X}_K(\xi_n), t_i)} \right]_v. \quad (2.77)$$

Inserting eqs. (2.75) and (2.76) into eq. (2.74) and introduction of the iteration index $r \in \{0, 1, \dots, R\} := \mathcal{R}$ leads to the following iterative scheme for the degrees of freedom of the cell-local space-time predictor:

$$\begin{aligned} & \underbrace{\int_K \left[\left[\hat{q}^{K,i,r+1} \right]_{nlv} \theta_{nlv}^{Ki} \Big|_{t_i+\Delta t_i} \right]_j \left[\theta_{\alpha\beta\gamma}^{Ki} \Big|_{t_i+\Delta t_i} \right]_j dx}_{\text{S-I}} - \\ & \underbrace{\int_{t_i}^{t_i+\Delta t_i} \int_K \left[\left[\hat{q}^{K,i,r+1} \right]_{nlv} \theta_{nlv}^{Ki} \right]_j \frac{\partial}{\partial t} \left[\theta_{\alpha\beta\gamma}^{Ki} \right]_j dx dt}_{\text{S-II}} = \\ & \underbrace{\int_K \left[\left[\hat{u}^{K,i} \right]_{nv} \phi_{nv}^K \right]_j \left[\theta_{\alpha\beta\gamma}^{Ki} \Big|_{t_i} \right]_j dx}_{\text{S-III}} + \\ & \underbrace{\int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \theta_{nlv}^{Ki} \right) \right]_{jk} \frac{\partial}{\partial x_k} \left[\theta_{\alpha\beta\gamma}^{Ki} \right]_j dx dt}_{\text{S-IV}} + \\ & \underbrace{\int_{t_i}^{t_i+\Delta t_i} \int_K \left[s \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \theta_{nlv}^{Ki} \right) \right]_j \left[\theta_{\alpha\beta\gamma}^{Ki} \right]_j dx dt}_{\text{S-V}}. \end{aligned} \quad (2.78)$$

We require this relation to hold for all $\alpha \in \mathcal{N}$, $\beta \in \mathcal{N}$ and $\gamma \in \mathcal{V}$.

As initial condition, i.e. for $r = 0$, we use

$$\left[\hat{q}^{K,i,0} \right]_{nvl} = \left[\hat{u}^{K,i} \right]_{nv} \quad (2.79)$$

for all time degrees of freedom $l \in \mathcal{N}$.

We will now proceed in a term-by-term fashion to rewrite all integrals with respect to reference coordinates so that we can finally derive a complete scheme to compute $\hat{\mathbf{q}}^{K,i,r+1}$ that holds for all $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

Term S-I

The first term of eq. (2.78) can be rewritten with respect to reference coordinates as follows:

$$\begin{aligned}
 & \int_K \left[\left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{nlw} \boldsymbol{\theta}_{nlv}^{Ki} \Big|_{t_i+\Delta t_i} \right]_j \left[\boldsymbol{\theta}_{\alpha\beta\gamma}^{Ki} \right]_{t_i+\Delta t_i} d\mathbf{x} = \\
 & \int_K \left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{nlv} \phi_n^K \left(\psi_l^i \Big|_{t_i+\Delta t_i} \right) [e_v]_j \phi_\alpha^K \left(\psi_\beta^i \Big|_{t_i+\Delta t_i} \right) [e_\gamma]_j d\mathbf{x} = \\
 & J_{\mathcal{X}_K} \int_{\hat{K}} \left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{nlv} \hat{\phi}_n \left(\hat{\psi}_l \Big|_1 \right) [e_v]_j \hat{\phi}_\alpha \left(\hat{\psi}_\beta \Big|_1 \right) [e_\gamma]_j d\boldsymbol{\xi} = \\
 & J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{nlv} \hat{\phi}_n(\hat{\boldsymbol{\xi}}_{\alpha'}) \left(\hat{\psi}_l \Big|_1 \right) [e_v]_j \hat{\phi}_\alpha(\hat{\boldsymbol{\xi}}_{\alpha'}) \left(\hat{\psi}_\beta \Big|_1 \right) [e_\gamma]_j \right) = \\
 & J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{nlv} \delta_{n\alpha'} \left(\hat{\psi}_l \Big|_1 \right) \delta_{vj} \delta_{\alpha\alpha'} \left(\hat{\psi}_\beta \Big|_1 \right) \delta_{j\gamma} \right) = \\
 & J_{\mathcal{X}_K} \hat{\omega}_\alpha \left[\hat{\psi}_\beta \Big|_1 \hat{\psi}_l \Big|_1 \right] \left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{\alpha l \gamma} = \\
 & J_{\mathcal{X}_K} \hat{\omega}_\alpha [\mathbf{R}]_{\beta,l} \left[\hat{\mathbf{q}}^{K,i,r+1} \right]_{\alpha l \gamma},
 \end{aligned} \tag{2.80}$$

where we remember from eq. (2.48) that

$$J_{\mathcal{X}_K} = \prod_{d=1}^D [\Delta \mathbf{x}]_d \tag{2.81}$$

and we define the matrix \mathbf{R} representing the Right Reference Element Mass Operator as

$$[\mathbf{R}]_{i,j} := \left[\hat{\psi}_i \Big|_1 \hat{\psi}_j \Big|_1 \right]_{i,j} \tag{2.82}$$

for $i, j \in \mathcal{N}$. A Python script to compute \mathbf{R} can be found in appendix A.

Term S-II

The second term of eq. (2.78) can be rewritten with respect to reference coordinates as follows:

$$\begin{aligned}
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[\hat{q}^{K,i,r+1} \right]_{nlv} \theta_{nlv}^{Ki} \left[\frac{\partial}{\partial t} \left[\theta_{\alpha\beta\gamma}^{Ki} \right]_j \right] dxdt = \\
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[\hat{q}^{K,i,r+1} \right]_{nlv} \phi_n^K \psi_l^i [e_v]_j \phi_\alpha^K \left(\frac{\partial}{\partial t} \psi_\beta^i \right) [e_\gamma]_j dxdt = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \int_0^1 \int_{\hat{K}} \left[\hat{q}^{K,i,r+1} \right]_{nlv} \hat{\phi}_n \hat{\psi}_l [e_v]_j \hat{\phi}_\alpha \left(\frac{1}{\Delta t_i} \frac{\partial}{\partial \tau} \hat{\psi}_\beta \right) [e_\gamma]_j d\xi d\tau = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[\hat{q}^{K,i,r+1} \right]_{nlv} \hat{\phi}_n(\hat{\xi}_{\alpha'}) \hat{\psi}_l(\hat{\tau}_{\beta'}) [e_v]_j \dots \right. \\
 & \quad \left. \dots \hat{\phi}_\alpha(\hat{\xi}_{\alpha'}) \left(\frac{\partial}{\partial \tau} \hat{\psi}_\beta(\hat{\tau}_{\beta'}) \right) [e_\gamma]_j \right) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[\hat{q}^{K,i,r+1} \right]_{nlv} \delta_{n\alpha'} \delta_{l\beta'} \delta_{vj} \dots \right. \\
 & \quad \left. \dots \delta_{\alpha\alpha'} \left(\frac{1}{\Delta t_i} \frac{\partial}{\partial \tau} \hat{\psi}_\beta(\hat{\tau}_{\beta'}) \right) \delta_{\gamma j} \right) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\alpha \frac{1}{\Delta t_i} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\beta'} \left[\frac{\partial}{\partial \tau} \hat{\psi}_\beta(\hat{\tau}_{\beta'}) \right] \left[\hat{q}^{K,i,r+1} \right]_{\alpha\beta'\gamma} \right) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\alpha \frac{1}{\Delta t_i} [\mathbf{K}]_{\beta,\beta'} \left[\hat{q}^{K,i,r+1} \right]_{\alpha\beta'\gamma} \tag{2.83}
 \end{aligned}$$

We recall from eq. (2.50) that

$$J_{\mathcal{T}_i} = \Delta t_i, \tag{2.84}$$

so that Δt_i and $1/\Delta t_i$ in eq. (2.83) cancel. In the derivation we made use of the fact that due to the chain rule

$$\frac{\partial}{\partial t} \psi_\beta^i = \frac{\partial}{\partial t} (\hat{\psi}_\beta \circ \mathcal{T}_i^{-1}) = \left(\frac{\partial}{\partial \tau} \hat{\psi}_\beta \right) \left(\frac{\partial}{\partial t} \mathcal{T}_i^{-1} \right) = \frac{1}{\Delta t_i} \frac{\partial}{\partial \tau} \hat{\psi}_\beta. \tag{2.85}$$

We furthermore introduce the matrix \mathbf{K} representing the Reference Element Stiffness Operator given as

$$[\mathbf{K}]_{ij} = \hat{\omega}_j \frac{\partial}{\partial \tau} \hat{\psi}_i(\hat{\tau}_j) \tag{2.86}$$

for $i, j \in \mathcal{N}$. A Python script to compute \mathbf{K} can be found in appendix A.

Term S-III

The third term of eq. (2.78) can be rewritten with respect to reference coordinates as

$$\begin{aligned}
 & \int_K \left[\left[\hat{\mathbf{u}}^{K,i} \right]_{nv} \boldsymbol{\phi}_{nv}^K \right]_j \left[\boldsymbol{\theta}_{\alpha\beta\gamma}^{Ki} \Big|_{t_i} \right]_j d\mathbf{x} = \\
 & \int_K \left[\hat{\mathbf{u}}^{K,i} \right]_{nv} \phi_n^K [e_v]_j \phi_\alpha^K \left(\psi_\beta^i \Big|_{t_i} \right) [e_\gamma]_j d\mathbf{x} = \\
 & J_{\mathcal{X}_K} \int_{\hat{K}} \left[\hat{\mathbf{u}}^{K,i} \right]_{nv} \hat{\phi}_n [e_v]_j \hat{\phi}_\alpha \left(\hat{\psi}_\beta \Big|_0 \right) [e_\gamma]_j d\boldsymbol{\xi} = \\
 & J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \left[\hat{\mathbf{u}}^{K,i} \right]_{nv} \hat{\phi}_n(\boldsymbol{\xi}_{\alpha'}) [e_v]_j \hat{\phi}_\alpha(\boldsymbol{\xi}_{\alpha'}) \left(\hat{\psi}_\beta \Big|_0 \right) [e_\gamma]_j \right) = \\
 & J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \left[\hat{\mathbf{u}}^{K,i} \right]_{nv} \delta_{n\alpha'} \delta_{vj} \delta_{\alpha\alpha'} \left(\hat{\psi}_\beta \Big|_0 \right) \delta_{\gamma j} \right) = \\
 & J_{\mathcal{X}_K} \hat{\omega}_\alpha \left[\hat{\psi}_\beta \Big|_0 \right] \left[\hat{\mathbf{u}}^{K,i} \right]_{\alpha\gamma} = \\
 & J_{\mathcal{X}_K} \hat{\omega}_\alpha [\mathbf{l}]_\beta \left[\hat{\mathbf{u}}^{K,i} \right]_{\alpha\gamma}, \tag{2.87}
 \end{aligned}$$

where we define the vector \mathbf{l} representing the Left Reference Element Flux Operator as

$$[\mathbf{l}]_i = \hat{\psi}_i \Big|_0 \tag{2.88}$$

for $i \in \mathcal{N}$. A Python script to compute \mathbf{l} can be found in appendix A.

Term S-IV

The third term of eq. (2.78) can be rewritten with respect to reference coordinates as

$$\begin{aligned}
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \theta_{nlv}^{Ki} \right) \right]_{jk} \frac{\partial}{\partial x_k} \left[\theta_{\alpha\beta\gamma}^{Ki} \right]_j dxdt = \\
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \phi_n^K \psi_l^i e_v \right) \right]_{jk} \left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([x]_d) \right) \psi_\beta^i(t) [e_\gamma]_j \dots \\
 & \dots \left(\frac{\partial}{\partial x_k} \psi_{[\alpha]_k}^K \right) dxdt = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \int_0^1 \int_{\hat{K}} \left[F \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \hat{\phi}_n \hat{\psi}_l e_v \right) \right]_{jk} \left(\prod_{d=1, d \neq k}^D \hat{\psi}_{[\alpha]_d}([\xi]_d) \right) \hat{\psi}_\beta(t) [e_\gamma]_j \dots \\
 & \dots \left(\frac{1}{[\Delta x]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\xi]_k) \right) d\xi d\tau = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[F \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \hat{\phi}_n(\hat{\xi}_{\alpha'}) \hat{\psi}_l(\hat{\tau}_{\beta'}) e_v \right) \right]_{jk} \dots \right. \\
 & \dots \left(\prod_{d=1, d \neq k}^D \hat{\psi}_{[\alpha]_d}([\hat{\xi}_{\alpha'}]_d) \right) \hat{\psi}_\beta(\hat{\tau}_{\beta'}) [e_\gamma]_j \left(\frac{1}{[\Delta x]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\hat{\xi}_{\alpha'}]_k) \right) \Bigg) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[F \left(\left[\hat{q}^{K,i,r} \right]_{nlv} \delta_{n\alpha'} \delta_{l\beta'} e_v \right) \right]_{jk} \dots \right. \\
 & \dots \left(\prod_{d=1, d \neq k}^D \delta_{[\alpha]_d}[\alpha']_d \right) \delta_{\beta\beta'} \delta_{\gamma j} \left(\frac{1}{[\Delta x]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\hat{\xi}_{\alpha'}]_k) \right) \Bigg) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\beta \sum_{k=1}^D \left(\frac{1}{[\Delta x]_k} \sum_{\alpha'_k \in \{0,1,\dots,N\}} \left(\prod_{d=0, d \neq k}^D \hat{\omega}_{[\alpha]_d} \dots \right. \right. \\
 & \dots \hat{\omega}_{\alpha'_k} \left(\frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}(\hat{\xi}_{\alpha'_k}) \right) \left[F \left(\left[\hat{q}^{K,i,r} \right]_{[\alpha]_{1:k-1}, \alpha'_k, [\alpha]_{k+1:N}} \right) \right]_{jk} \Bigg) \Bigg) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\beta \sum_{k=1}^D \left(\frac{1}{[\Delta x]_k} \sum_{\alpha'_k \in \{0,1,\dots,N\}} \left(\prod_{d=0, d \neq k}^D \hat{\omega}_{[\alpha]_d} \dots \right. \right. \\
 & \dots [K]_{[\alpha]_k, \alpha'_k} \left[F \left(\left[\hat{q}^{K,i,r} \right]_{[\alpha_0, \alpha_1, \dots, \alpha_{k-1}, \alpha'_k, \alpha_{k+1}, \dots, \alpha_N] \beta v} e_v \right) \right]_{\gamma k} \Bigg) \Bigg),
 \end{aligned}$$

where we used that

$$\begin{aligned}
 \frac{\partial}{\partial x_k} \theta_{\alpha\beta\gamma}^{Ki}(\mathbf{x}, t) &= \left(\frac{\partial}{\partial x_k} \phi_{\alpha}^K(\mathbf{x}) \right) \psi_{\beta}^i(t) \mathbf{e}_{\gamma} = \left(\frac{\partial}{\partial x_k} \prod_{d=1}^D \psi_{[\alpha]_d}^K([\mathbf{x}]_d) \right) \psi_{\beta}^i(t) \mathbf{e}_{\gamma} = \\
 &= \left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([\mathbf{x}]_d) \right) \left(\frac{\partial}{\partial x_k} \psi_{[\alpha]_k}^K([\mathbf{x}]_k) \right) \psi_{\beta}^i(t) \mathbf{e}_{\gamma} = \\
 &= \left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([\mathbf{x}]_d) \right) \left(\frac{\partial}{\partial x_k} \hat{\psi}_{[\alpha]_k} \left([\boldsymbol{\chi}_K^{-1}(\mathbf{x})]_k \right) \right) \psi_{\beta}^i(t) \mathbf{e}_{\gamma} = \\
 &= \left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([\mathbf{x}]_d) \right) \left(\left(\frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k} \left([\boldsymbol{\chi}_K^{-1}(\mathbf{x})]_k \right) \right) \left(\frac{\partial}{\partial x_k} [\boldsymbol{\chi}_K^{-1}(\mathbf{x})]_k \right) \right) \dots \\
 &\dots \psi_{\beta}^i(t) \mathbf{e}_{\gamma} = \\
 &= \left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([\mathbf{x}]_d) \right) \left(\frac{1}{[\Delta \mathbf{x}^K]_k} \frac{\partial}{\partial \xi_k} \hat{\phi}_{[\alpha]_k} \left([\boldsymbol{\chi}_K^{-1}(\mathbf{x})]_k \right) \right) \psi_{\beta}^i(t) \mathbf{e}_{\gamma}.
 \end{aligned} \tag{2.90}$$

Term S-V

The fifth term of eq. (2.78) can be rewritten with respect to reference coordinates as follows:

$$\begin{aligned}
 &\int_{t_i}^{t_i + \Delta t_i} \int_K \left[\mathbf{s} \left([\hat{\mathbf{q}}^{K,i,r}]_{nlv} \theta_{nlv}^{Ki} \right) \right]_j [\theta_{\alpha\beta\gamma}^{Ki}]_j d\mathbf{x} dt = \\
 &J_{\mathcal{T}_i} J_{\mathcal{X}_K} \int_0^1 \int_{\hat{K}} \left[\mathbf{s} \left([\hat{\mathbf{q}}^{K,i,r}]_{nlv} \hat{\phi}_n \hat{\psi}_l \mathbf{e}_v \right) \right]_j \hat{\phi}_{\alpha} \hat{\psi}_l [\mathbf{e}_{\gamma}]_j d\xi d\tau = \\
 &J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[\mathbf{s} \left([\hat{\mathbf{q}}^{K,i,r}]_{nlv} \hat{\phi}_n(\xi_{\alpha'}) \hat{\psi}_l(\hat{\tau}_{\beta'}) \mathbf{e}_v \right) \right]_j \dots \right. \\
 &\dots \left. \hat{\phi}_{\alpha}(\xi_{\alpha'}) \hat{\psi}_{\beta}(\hat{\tau}_{\beta'}) [\mathbf{e}_{\gamma}]_j \right) = \\
 &J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[\mathbf{s} \left([\hat{\mathbf{q}}^{K,i,r}]_{nlv} \delta_{n\alpha'} \delta_{l\beta'} \mathbf{e}_v \right) \right]_j \delta_{\alpha\alpha'} \delta_{\beta\beta'} \delta_{\gamma j} \right) = \\
 &J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_{\alpha} \hat{\omega}_{\beta} \left[\mathbf{s} \left([\hat{\mathbf{q}}^{K,i,r}]_{\alpha\beta v} \mathbf{e}_v \right) \right]_{\gamma}.
 \end{aligned} \tag{2.91}$$

The Complete Fixed-point Iteration

Now collecting the results from eqs. (2.80), (2.83), (2.87), (2.89) and (2.91), plugging them back into eq. (2.78) and division of the resulting equation by $J_{\mathcal{X}_K} \hat{\omega}_\alpha$ yields

$$\begin{aligned}
 (\mathbf{R} - \mathbf{K})_{\beta\beta'} [\hat{\mathbf{q}}^{K,i,r+1}]_{\alpha\beta'\gamma} &= [\mathbf{I}]_\beta [\hat{\mathbf{u}}^{K,i}]_{\alpha\gamma} + \\
 J_{\mathcal{T}_i} \frac{\hat{\omega}_\beta}{\hat{\omega}_\alpha} \sum_{k=1}^D \left(\frac{1}{[\Delta \mathbf{x}]_k} \sum_{\alpha'_k \in \{0,1,\dots,N\}} \left(\prod_{d=0, d \neq k}^D \hat{\omega}_{[\alpha]_d} \dots \right. \right. \\
 \left. \left. \dots [\mathbf{K}]_{[\alpha]_k, \alpha'_k} \left[\mathbf{F} \left([\hat{\mathbf{q}}^{K,i,r}]_{[\alpha_0, \alpha_1, \dots, \alpha_{k-1}, \alpha'_k, \alpha_{k+1}, \dots, \alpha_N] \beta v} \mathbf{e}_v \right) \right]_{\gamma k} \right) \right) + \\
 J_{\mathcal{T}_i} \hat{\omega}_\beta \left[\mathbf{s} \left([\hat{\mathbf{q}}^{K,i,r}]_{\alpha\beta v} \mathbf{e}_v \right) \right]_\gamma, \tag{2.92}
 \end{aligned}$$

which has to hold for all $\alpha \in \mathcal{N}$, $\beta \in \mathcal{N}$ and $\gamma \in \mathcal{V}$. To speed up the computation of the new iterate $\hat{\mathbf{q}}^{K,i,r+1}$ we can invert the matrix on the left-hand side prior to the simulation to obtain the iteration matrix $\tilde{\mathbf{K}} := (\mathbf{R} - \mathbf{K})^{-1}$. A Python script to compute $\tilde{\mathbf{K}}$ can be found in appendix A.

A possible termination criterion could be $\Delta < \varepsilon$, where $\varepsilon > 0$ is a suitable constant related to the desired accuracy of the iteration, e.g. $\varepsilon = 10^{-7}$ and the squared element-wise residual Δ^2 is defined as follows:

$$\Delta^2 = \sum_{n \in \mathcal{N}} \sum_{l \in \mathcal{N}} \sum_{v \in \mathcal{V}} \left([\hat{\mathbf{q}}^{K,i,r+1}]_{n,l,v} - [\hat{\mathbf{q}}^{K,i,r}]_{n,l,v} \right)^2. \tag{2.93}$$

For linear homogeneous scalar hyperbolic balance laws and neglecting floating point errors it can be proven that the iteration converges after at most N steps (see [10] for details).

2.1.12 A Fully-discrete Update Scheme for the Time-discrete Solution

Now that we have developed a method to compute the space-time predictor, we can go back to the original one-step, cell-local update scheme given in eq. (2.34). Inserting the local space-time predictor $\tilde{q}_h^{K,i}$ yields

$$\begin{aligned} \int_K \left[\tilde{u}_h^{K,i} \Big|_{t_i+\Delta t_i} \right]_v \left[w_h^K \right]_v dx &= \int_K \left[\tilde{u}_h^{K,i} \Big|_{t_i} \right]_v \left[w_h^K \right]_v dx + \\ &\int_{t_i}^{t_i+\Delta t_i} \int_K \left[F(\tilde{q}_h^{K,i}) \right]_{vd} \frac{\partial}{\partial x_d} \left[w_h^K \right]_v dx dt + \\ &\int_{t_i}^{t_i+\Delta t_i} \int_K \left[s(\tilde{q}_h^{K,i}) \right]_v \left[w_h^K \right]_v dx dt - \\ &\int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} \left[\mathcal{G}(\tilde{q}_h^{K,i}, \tilde{q}_h^{K+i}, n) \right]_v \left[w_h^K \right]_v ds(x) dt, \end{aligned} \quad (2.94)$$

which has to hold for all $v \in \mathcal{V}$, $K \in \mathcal{K}_h$, $w_h \in \mathbb{W}_h$ and $i \in \mathcal{I}$.

Making use of the bases we derived earlier the cell-local solution $\tilde{u}_h^{K,i}$ at times t_i and $t_i + \Delta t_i$ can be represented by tensors of coefficients $\hat{u}^{K,i}$ and $\hat{u}^{K,i+1}$ as

$$\tilde{u}_h^{K,i} \Big|_{t_i} = \left[\hat{u}^{K,i} \right]_{n,v} \phi_{n,v}^K \quad (2.95)$$

and

$$\tilde{u}_h^{K,i} \Big|_{t_i+\Delta t_i} = \left[\hat{u}^{K,i+1} \right]_{n,v} \phi_{n,v}^K, \quad (2.96)$$

respectively. Inserting eqs. (2.95) and (2.96) and the ansatz for the space-time predictor (2.75) into eq. (2.94) yields

$$\begin{aligned} &\underbrace{\int_K \left[\left[\hat{u}^{K,i+1} \right]_{n,v} \phi_{n,v}^K \right]_j \left[\phi_{\alpha,\gamma}^K \right]_j dx}_{\text{U-I}} = \underbrace{\int_K \left[\left[\hat{u}^{K,i} \right]_{n,v} \phi_{n,v}^K \right]_j \left[\phi_{\alpha,\gamma}^K \right]_j dx}_{\text{U-II}} + \\ &\underbrace{\int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left(\left[\hat{q}^{K,i} \right]_{n,l,v} \theta_{n,l,v}^{Ki} \right) \right]_{jk} \frac{\partial}{\partial x_k} \left[\phi_{\alpha,\gamma}^K \right]_j dx dt}_{\text{U-III}} + \\ &\underbrace{\int_{t_i}^{t_i+\Delta t_i} \int_K \left[s \left(\left[\hat{q}^{K,i} \right]_{n,l,v} \theta_{n,l,v}^{Ki} \right) \right]_j \left[\phi_{\alpha,\gamma}^K \right]_j dx dt}_{\text{U-IV}} - \\ &\underbrace{\int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+i}, n \right) \right]_j \left[\phi_{\alpha,\gamma}^K \right]_j ds(x) dt}_{\text{U-V}}, \end{aligned} \quad (2.97)$$

2.1. A D-dimensional ADER-DG Scheme With MUSCL-Hancock a Posteriori Subcell Limiting for Non-linear Hyperbolic Balance Laws

which we require to hold for all $\alpha \in \mathcal{N}$, $\gamma \in \mathcal{V}$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$. In the following we will again proceed by simplifying each term in reference coordinates separately and then in the end assemble all terms to obtain a complete fully-discrete update scheme.

Term U-I

The first term of eq. (2.97) can be rewritten with respect to reference coordinates as follows:

$$\begin{aligned}
 & \int_K \left[\left[\hat{\mathbf{u}}^{K,i+1} \right]_{n,v} \boldsymbol{\phi}_{n,v}^K \right]_j \left[\boldsymbol{\phi}_{\alpha,\gamma}^K \right]_j dx = \\
 & \int_K \left[\left[\hat{\mathbf{u}}^{K,i+1} \right]_{n,v} \phi_n^K \mathbf{e}_v \right]_j \left[\phi_\alpha^K \mathbf{e}_\gamma \right]_j dx = \\
 & J_{\mathcal{X}_K} \int_{\hat{K}} \left[\left[\hat{\mathbf{u}}^{K,i+1} \right]_{n,v} \hat{\phi}_n \mathbf{e}_v \right]_j \left[\hat{\phi}_\alpha \mathbf{e}_\gamma \right]_j d\hat{\xi} = \\
 & J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \left[\hat{\mathbf{u}}^{K,i+1} \right]_{n,v} \hat{\phi}_n(\hat{\xi}_{\alpha'}) [\mathbf{e}_v]_j \hat{\phi}_\alpha(\hat{\xi}_{\alpha'}) [\mathbf{e}_\gamma]_j \right) = \\
 & J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \left[\hat{\mathbf{u}}^{K,i+1} \right]_{n,v} \delta_{n\alpha'} \delta_{vj} \delta_{\alpha\alpha'} \delta_{\gamma j} \right) = \\
 & J_{\mathcal{X}_K} \hat{\omega}_\alpha \left[\hat{\mathbf{u}}^{K,i+1} \right]_{\alpha,\gamma}. \tag{2.98}
 \end{aligned}$$

Term U-II

Analogously to the first term of eq. (2.97), the second term can be rewritten as follows:

$$\begin{aligned}
 & \int_K \left[\left[\hat{\mathbf{u}}^{K,i} \right]_{n,v} \boldsymbol{\phi}_{n,v}^K \right]_j \left[\boldsymbol{\phi}_{\alpha,\gamma}^K \right]_j dx = \\
 & J_{\mathcal{X}_K} \hat{\omega}_\alpha \left[\hat{\mathbf{u}}^{K,i} \right]_{\alpha,\gamma}. \tag{2.99}
 \end{aligned}$$

Term U-III

The third term of eq. (2.97) can be rewritten with respect to reference coordinates as

$$\begin{aligned}
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left([\hat{q}^{K,i}]_{n,l,v} \theta_{n,l,v}^{Ki} \right) \right]_{jk} \frac{\partial}{\partial x_k} [\phi_{\alpha,\gamma}^K]_j dx dt = \\
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left([\hat{q}^{K,i}]_{n,l,v} \phi_n^K \psi_l^i e_v \right) \right]_{jk} \frac{\partial}{\partial x_k} \left(\prod_{d=1}^D \psi_{[\alpha]_d}^K([x]_d) \right) [e_\gamma]_j dx dt = \\
 & \int_{t_i}^{t_i+\Delta t_i} \int_K \left[F \left([\hat{q}^{K,i}]_{n,l,v} \phi_n^K \psi_l^i e_v \right) \right]_{jk} \left(\prod_{d=1,d \neq k}^D \psi_{[\alpha]_d}^K([x]_d) \right) \frac{1}{[\Delta x^K]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\mathbf{x}_K(x)]_k) [e_\gamma]_j dx dt = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \int_0^1 \int_{\hat{K}} \left[F \left([\hat{q}^{K,i}]_{n,l,v} \hat{\phi}_n \hat{\psi}_l e_v \right) \right]_{kj} \left(\prod_{d=1,d \neq k}^D \hat{\psi}_{[\alpha]_d}([\xi]_d) \right) \frac{1}{[\Delta \mathbf{x}^K]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\xi]_k) [e_\gamma]_j d\xi d\tau = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[F \left([\hat{q}^{K,i}]_{n,l,v} \hat{\phi}_n(\xi_{\alpha'}) \hat{\psi}(\hat{\tau}_{\beta'}) e_v \right) \right]_{jk} \left(\prod_{d=1,d \neq k}^D \hat{\psi}_{[\alpha]_d}([\xi_{\alpha'}]_d) \right) \frac{1}{[\Delta \mathbf{x}^K]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\xi_{\alpha'}]_k) \right) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[F \left([\hat{q}^{K,i}]_{n,l,v} \delta_{n\alpha'} \delta_{l\beta'} e_v \right) \right]_{jk} \left(\prod_{d=1,d \neq k}^D \delta_{[\alpha]_d}[\alpha']_d \right) \frac{1}{[\Delta \mathbf{x}^K]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\xi_{\alpha'}]_k) \right) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\alpha \sum_{k=1}^D \left(\sum_{\alpha'_k \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\frac{\hat{\omega}_{\beta'}}{\hat{\omega}_{\alpha'_k}} \frac{1}{[\Delta \mathbf{x}^K]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{\alpha'_k}([\xi]_{\alpha'_k}) \right) \left[F \left([\hat{q}^{K,i}]_{[\alpha]_1, [\alpha]_2, \dots, [\alpha]_{k-1}, \alpha'_k, [\alpha]_{k+1}, \dots, [\alpha]_D} \right) \right]_{\beta', v} e_v \right) = \\
 & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\alpha \sum_{k=1}^D \left(\sum_{\alpha'_k \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\frac{1}{\hat{\omega}_{\alpha'_k}} \frac{1}{[\Delta \mathbf{x}^K]_k} [\mathbf{K}]_{\alpha'_k, k} \left[F \left([\hat{q}^{K,i}]_{[\alpha]_1, [\alpha]_2, \dots, [\alpha]_{k-1}, \alpha'_k, [\alpha]_{k+1}, \dots, [\alpha]_D} \right) \right]_{\beta', v} e_v \right) \right)_{\gamma, k}
 \end{aligned} \tag{2.100}$$

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where we made use of the fact that due to the chain rule it holds that

$$\begin{aligned}
\frac{\partial}{\partial x_k} \left(\prod_{d=1}^D \psi_{[\alpha]_d}^K([x]_d) \right) &= \left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([x]_d) \right) \frac{\partial}{\partial x_k} \psi_{[\alpha]_k}^K([x]_k) = \\
&\left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([x]_d) \right) \frac{\partial}{\partial \xi_j} \hat{\psi}_{[\alpha]_k}([\mathbf{x}_K(\mathbf{x})]_k) \frac{\partial}{\partial x_k} [\mathbf{x}_K(\mathbf{x})]_j = \\
&\left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([x]_d) \right) \frac{\partial}{\partial \xi_j} \hat{\psi}_{[\alpha]_k}([\mathbf{x}_K(\mathbf{x})]_k) \frac{1}{[\Delta \mathbf{x}^K]_k} \delta_{kj} = \\
&\left(\prod_{d=1, d \neq k}^D \psi_{[\alpha]_d}^K([x]_d) \right) \frac{1}{[\Delta \mathbf{x}^K]_k} \frac{\partial}{\partial \xi_k} \hat{\psi}_{[\alpha]_k}([\mathbf{x}_K(\mathbf{x})]_k) dx dt. \tag{2.101}
\end{aligned}$$

Term U-IV

The fourth term of eq. (2.97) can be rewritten with respect to reference coordinates as follows:

$$\begin{aligned}
&\int_{t_i}^{t_i+\Delta t_i} \int_K \left[s \left([\hat{\mathbf{q}}^{K,i}]_{n,l,v} \boldsymbol{\theta}_{n,l,v}^{Ki} \right) \right]_j [\boldsymbol{\phi}_{\alpha,\gamma}^K]_j dx dt = \\
&\int_{t_i}^{t_i+\Delta t_i} \int_K \left[s \left([\hat{\mathbf{q}}^{K,i}]_{n,l,v} \boldsymbol{\phi}_n^K \psi_l^i \mathbf{e}_v \right) \right]_j \boldsymbol{\phi}_\alpha^K [\mathbf{e}_\gamma]_j dx dt = \\
&J_{\mathcal{T}_i} J_{\mathcal{X}_K} \int_0^1 \int_{\hat{K}} \left[s \left([\hat{\mathbf{q}}^{K,i}]_{n,l,v} \hat{\boldsymbol{\phi}}_n \hat{\psi}_l \mathbf{e}_v \right) \right]_j \hat{\boldsymbol{\phi}}_\alpha [\mathbf{e}_\gamma]_j d\hat{\xi} d\tau = \\
&J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[s \left([\hat{\mathbf{q}}^{K,i}]_{n,l,v} \hat{\boldsymbol{\phi}}_n(\hat{\boldsymbol{\xi}}_{\alpha'}) \hat{\psi}_l(\hat{\tau}_{\beta'} \mathbf{e}_v) \right) \right]_j \hat{\boldsymbol{\phi}}_\alpha(\hat{\boldsymbol{\xi}}_{\alpha'}) [\mathbf{e}_\gamma]_j \right) = \\
&J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\alpha' \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\alpha'} \hat{\omega}_{\beta'} \left[s \left([\hat{\mathbf{q}}^{K,i}]_{n,l,v} \delta_{n\alpha'} \delta_{l\beta'} \mathbf{e}_v \right) \right]_j \delta_{\alpha\alpha'} \delta_{\gamma j} \right) = \\
&J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_\alpha \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\beta'} \left[s \left([\hat{\mathbf{q}}^{K,i}]_{\alpha,\beta',v} \mathbf{e}_v \right) \right]_\gamma \right). \tag{2.102}
\end{aligned}$$

Term U-V

Let $d \in \mathcal{D}$ and $e \in \{0,1\} := \mathcal{E}$. Then if we define the $D-1$ -dimensional quadrilateral $\partial \hat{K}_{d,e}$ as

$$\partial \hat{K}_{d,e} = \left\{ \boldsymbol{\xi} \in \hat{K} \mid [\boldsymbol{\xi}]_d = e \right\}, \tag{2.103}$$

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the set $\{\partial\hat{K}_{d,e}\}_{d\in\mathcal{D},e\in\mathcal{E}}$ is a partition of the surface $\partial\hat{K}$ of the spatial reference element. Utilizing the mappings \mathcal{X}_K that map points $\xi \in \hat{K}$ to $x \in K$ for all $K \in \mathcal{K}_h$ we can define

$$\partial K_{d,e} = \mathcal{X}_K \left(\partial\hat{K}_{d,e} \right), \quad (2.104)$$

where now the set $\{\partial K_{d,e}\}_{d\in\mathcal{D},e\in\mathcal{E}}$ is a quadrilateral partition of the surface ∂K for all cells $K \in \mathcal{K}_h$.

In consequence the surface integral in the fifth term of eq. (2.97) can be rewritten as follows:

$$\begin{aligned} & \int_{t_i}^{t_i+\Delta t_i} \int_{\partial K} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, n \right) \right]_j \left[\phi_{\alpha,\gamma}^K \right]_j ds(x) dt = \\ & \int_{t_i}^{t_i+\Delta t_i} \sum_{d\in\mathcal{D}} \sum_{e\in\mathcal{E}} \left(\int_{\partial K_{d,e}} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, e_d \right) \right]_j \phi_{\alpha}^K \left[e_{\gamma} \right]_j ds(x) \right) dt = \\ & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \int_0^1 \sum_{d\in\mathcal{D}} \sum_{e\in\mathcal{E}} \left(\frac{1}{[\Delta x^K]_d} \int_{\partial\hat{K}_{d,e}} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, (-1)^e e_d \right) \right]_j \hat{\phi}_{\alpha} \left[e_d \right]_j ds(\xi) \right) d\tau = \\ & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\beta'\in\mathcal{D}} \hat{\omega}_{\beta'} \sum_{d\in\mathcal{D}} \sum_{e\in\mathcal{E}} \sum_{\alpha'\in\mathcal{N}^-} \left(\hat{\omega}_{\alpha'} \frac{1}{[\Delta x^K]_d} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, (-1)^e e_d \right) \right]_j \hat{\phi}_{\alpha'}(\hat{\xi}_{\alpha'}) \left(\hat{\psi}_{[\alpha]_d|_e} \right) \left[e_d \right]_j \right) = \\ & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \sum_{\beta'\in\mathcal{D}} \hat{\omega}_{\beta'} \sum_{d\in\mathcal{D}} \sum_{e\in\mathcal{E}} \sum_{\alpha'\in\mathcal{N}^-} \left(\hat{\omega}_{\alpha'} \frac{1}{[\Delta x^K]_d} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, (-1)^e e_d \right) \right]_j \delta_{\alpha'd} \left(\hat{\psi}_{[\alpha]_d|_e} \right) \delta_{\gamma j} \right) = \\ & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_{\alpha} \sum_{\beta'\in\mathcal{D}} \sum_{d\in\mathcal{D}} \sum_{e\in\mathcal{E}} \sum_{\alpha'_d\in\mathcal{N}} \left(\frac{\hat{\omega}_{\beta'}}{\hat{\omega}_{\alpha'_d}} \frac{1}{[\Delta x^K]_d} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, (-1)^e e_d \right) \right]_{\gamma} \left(\hat{\psi}_{\alpha'_d|_e} \right) \right) = \\ & J_{\mathcal{T}_i} J_{\mathcal{X}_K} \hat{\omega}_{\alpha} \sum_{\beta'\in\mathcal{D}} \sum_{d\in\mathcal{D}} \sum_{e\in\mathcal{E}} \sum_{\alpha'_d\in\mathcal{N}} \left(\frac{\hat{\omega}_{\beta'}}{\hat{\omega}_{\alpha'_d}} \frac{1}{[\Delta x^K]_d} \left[\mathcal{G} \left(\hat{q}^{K,i}, \hat{q}^{K+,i}, (-1)^e e_d \right) \right]_{\gamma} \left(\delta_{e0} \mathbf{l}_{\alpha'_d} \right) \left(\delta_{e1} \mathbf{r}_{\alpha'_d} \right) \left(\hat{\psi}_{\alpha'_d|_e} \right) \right). \end{aligned} \quad (2.105)$$

In each term we have to solve a Riemann problem in direction of the unit vector e_d defined as

$$[e_d]_{d'} = \delta_{dd'} \quad (2.106)$$

for $d' \in \mathcal{D}$. \mathbf{l} and \mathbf{r} denote the Left and Right Reference Element Flux Operator, respectively, and the latter is defined as

$$[\mathbf{r}]_i = \hat{\psi}_i \Big|_1 \quad (2.107)$$

for $i \in \mathcal{N}$. A Python script to compute \mathbf{r} can be found in appendix A.

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The Complete One-step Update Scheme

Inserting eqs. (2.98) to (2.100), (2.102) and (2.105) into eq. (2.97) and dividing the resulting equation by $\hat{\omega}_\alpha$ and $J_{\mathcal{X}_K}$ yields

$$\begin{aligned}
 [\hat{\mathbf{u}}^{K,i+1}]_{\alpha,\gamma} &= [\hat{\mathbf{u}}^{K,i}]_{\alpha,\gamma} + \\
 &J_{\mathcal{T}_i} \sum_{k=1}^D \left(\sum_{\alpha'_k \in \mathcal{N}} \sum_{\beta' \in \mathcal{N}} \left(\frac{\hat{\omega}_{\beta'}}{\hat{\omega}_{\alpha'_k}} \frac{1}{[\Delta \mathbf{x}^K]_k} \underbrace{\frac{\partial}{\partial \xi_k} \hat{\psi}_{\alpha'_k} \left([\hat{\xi}]_{\alpha'_k} \right)}_{\text{Kxi}_{\alpha'_k k}} \left[F \left([\hat{\mathbf{q}}^{K,i}]_{[\alpha]_1, [\alpha]_2, \dots, [\alpha]_{k-1}, \alpha'_k, [\alpha]_{k+1}, \dots, [\alpha]_D}, \beta', v \right) \right]_{\gamma, k} \right) \right) + \\
 &J_{\mathcal{T}_i} \sum_{\beta' \in \mathcal{N}} \left(\hat{\omega}_{\beta'} \left[s \left([\hat{\mathbf{q}}^{K,i}]_{\alpha, \beta', v} \right) \right]_{\gamma} \right) - \\
 &J_{\mathcal{T}_i} \sum_{\beta' \in \mathcal{D}} \sum_{d \in \mathcal{D}} \sum_{e \in \mathcal{E}} \sum_{\alpha'_d \in \mathcal{N}} \left(\frac{\hat{\omega}_{\beta'}}{\hat{\omega}_{\alpha'_d}} \frac{1}{[\Delta \mathbf{x}^K]_d} \left[\mathcal{G} \left(\hat{\mathbf{q}}^{K,i}, \hat{\mathbf{q}}^{K+,i}, (-1)^e \mathbf{e}_d \right) \right]_{\gamma} \underbrace{\left(\hat{\psi}_{\alpha'_d} \right)_e}_{\text{F0, F1}} \right), \tag{2.108}
 \end{aligned}$$

which we require to hold for $\alpha \in \mathcal{N}$, $\gamma \in \mathcal{V}$, $K \in \mathcal{K}_h$ and $i \in \mathcal{I}$.

Time step restriction

For the scheme to be stable we require that the following inequality holds for the time step Δt_i :

$$\Delta t_i \leq \frac{1}{D} \frac{1}{(2N+1)} \min_{d \in \mathcal{D}} \left(\frac{[\Delta \mathbf{x}]_d}{\Lambda^{i,d}} \right), \tag{2.109}$$

where

$$\Lambda^{i,d} = \max_{v \in \mathcal{V}} \text{abs} \left[\lambda^{i,d} \right]_v \tag{2.110}$$

and $\lambda^{d,i}$ is a vector containing the V real eigenvalues of the Jacobian

$$\frac{\partial}{\partial x_k} \left[F(\mathbf{u}(\mathbf{x}, t_i)) \right]_{jd} \tag{2.111}$$

for the respective dimension $d \in \mathcal{D}$ and the index of the current time step $i \in \mathcal{I}$. For details on the derivation of this formula see [11, 3, 12].

2.1.13 A Posteriori Subcell Limiting

The unlimited ADER-DG scheme derived in the previous section allows us to solve non-linear hyperbolic balance laws with arbitrary order in both time and space for continuous data. For discontinuous initial data and in scenarios where even for smooth initial data due to the non-linear nature of the system shocks arise, however, our high-order discontinuous Galerkin method is unsuitable. It is a linear scheme in the sense of Godunov, i.e. the coefficients of the scheme are independent of the current state of the system (see [13, 2] for a derivation of the so-called Godunov theorem) and therefore shocks in the system do not only give rise to artificial and persistent oscillations, but also decrease pointwise accuracy in the vicinity to first order and even cause loss of pointwise convergence at the point of discontinuity (see [6] for details). The quality of the numerical solution is unacceptable at best, if not the simulation crashes altogether due to a strictly positive physical quantity such as pressure or density becoming negative and thereby invalidating the well-posedness of the problem.

A way to solving this issue is limiting. For an exhaustive overview on classical limiting approaches see again [2]. According to Dumbser et al. [3], there are two key challenges in designing limiter procedures:

1. A so-called troubled cell indicator needs to implement criteria on how to identify troubled cells, i.e. cells that need limiting.
2. The troubled DG solution needs to be replaced by a robust non-linear reconstructions scheme in a way such that the minimum amount of artificial numerical viscosity is injected in these cells ideally preserving the subcell resolution property of the high-order DG scheme.

In the following we will present a novel approach called a posterior subcell limiting. It was first introduced in 2014 (see [3]) and has very favorable properties with respect to the aforementioned challenges. The general idea is to first project the local ADER-DG solution onto a finer equidistant grid to check if the solution satisfies certain admissibility criteria. If this is not the case, we discard this candidate solution and instead rely on the more robust MUSCL-Hancock FVM scheme starting from the fine grid solution of the previous time step (which is either the projected ADER-DG solution or, if the cell had already been troubled in the previous time step, the available FVM fine grid solution). For troubled cells we finally use a reconstruction operator to replace the rejected candidate solution by the transformed fine grid solution. The remainder of this chapter is structured as follows: We will first introduce the necessary projection and reconstruction operators. Subsequently we will discuss a possible set of admissibility criteria and conclude with a summary on the MUSCL-Hancock finite volume method as

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an example for a robust scheme that can be used for a posteriori subcell limiting.

Projection and Reconstruction

In order to check for admissibility and in case of a troubled cell to employ a more robust FVM scheme, we need to project the ADER-DG degrees of freedom $\hat{\mathbf{u}}^{K,i}$ to N_S^D subcell averages $\hat{\mathbf{p}}^{K,i}$ on an equidistant fine grid. We choose to split $K \in \mathcal{K}_h$ into $N_S = 2N + 1$ subcells along each spatial dimension creating a total of N_S^D subcells denoted as K_α , $\alpha \in \{0, 1, \dots, 2N\} := \mathcal{N}_S$. For explicit Godunov-type finite volume schemes on the subgrid we must satisfy the stability condition

$$\Delta t \leq \frac{1}{d} \frac{1}{N_S} \min_{d \in \mathcal{D}} \left(\frac{[\Delta \mathbf{x}]_d}{\Lambda^d} \right). \quad (2.112)$$

Comparing eq. (2.112) to the time step restriction for the ADER-DG scheme given in eq. (2.109) illustrates that the choice $N_S = 2N + 1$ is optimal in the sense that it makes sure that, on the one hand, time steps on the ADER-DG grid are also stable on the equidistant subgrid and, on the other hand, that we add the minimum amount of dissipation necessary. See [3] for additional remarks on the optimality of this particular choice of N_S .

Let as before K_α be a cell in the equidistant subgrid of cardinality $N_S^D = (2N + 1)^D$ on a cell $K \in \mathcal{K}_h$. Then we can define a projection \mathcal{P} of the degrees of freedom from the ADER-DG grid cell K to the subcell K_α for all $\alpha \in \mathcal{N}_S$ by prescribing that the integral averages over K_α are to be preserved. Mathematically this requirement can be expressed as

$$\begin{aligned} [\hat{\mathbf{p}}^{K,i}]_{\alpha,\gamma} &= \frac{1}{|K_\alpha|} \int_{K_\alpha} [\hat{\mathbf{u}}_h^{K,i}]_\gamma dx = \frac{1}{|K_\alpha|} \int_{K_\alpha} [\hat{\mathbf{u}}^{K,i}]_{n,v} \phi_n^K(\mathbf{x}) [e_v]_\gamma dx = \\ &= \sum_{n \in \mathcal{N}} \sum_{\alpha' \in \mathcal{N}} \left([\hat{\mathbf{u}}^{K,i}]_{n,\gamma} \hat{\omega}_{\alpha'} \hat{\phi}_n \left(\frac{1}{N_S} \alpha + \frac{1}{N_S} \hat{\xi}_{\alpha'} \right) \right) = \\ &= \sum_{n \in \mathcal{N}} \sum_{\alpha' \in \mathcal{N}} \left(\prod_{d \in \mathcal{D}} \left(\hat{\omega}_{\alpha'_d} \hat{\psi}_{n_d} \left(\frac{1}{N_S} \alpha_d + \frac{1}{N_S} \hat{\xi}_{\alpha'_d} \right) \right) [\hat{\mathbf{u}}^{K,i}]_{n,\gamma} \right) = \\ &= \sum_{n \in \mathcal{N}} \left(\sum_{\alpha'_0 \in \mathcal{N}} \left(\hat{\omega}_{\alpha'_0} \hat{\psi}_{n_0} \left(\frac{1}{N_S} \alpha_0 + \frac{1}{N_S} \hat{\xi}_{\alpha'_0} \right) \dots \sum_{\alpha'_{D-1} \in \mathcal{N}} \left(\hat{\omega}_{\alpha'_{D-1}} \hat{\psi}_{n_{D-1}} \left(\frac{1}{N_S} \alpha_{D-1} + \frac{1}{N_S} \hat{\xi}_{\alpha'_{D-1}} \right) \right) \right) [\hat{\mathbf{u}}^{K,i}]_n \right) \\ &= \sum_{n \in \mathcal{N}} \left(\prod_{d \in \mathcal{D}} ([P]_{\alpha_d, n_d}) [\hat{\mathbf{u}}^{K,i}]_{n,\gamma} \right), \end{aligned} \quad (2.113)$$

where $|K_\alpha|$ denotes the volume of K_α for $\alpha \in \mathcal{N}$ and $\gamma \in \mathcal{D}$. The computation can be carried out efficiently in a dimension-by-dimension manner using the projection matrix

$$[P]_{ij} = \sum_{k \in \mathcal{N}} \left(\hat{\omega}_k \hat{\psi}_j \left(\frac{1}{N_S} i + \frac{1}{N_S} \hat{\xi}_k \right) \right) \quad (2.114)$$

for $i \in \mathcal{N}_S$ and $j \in \mathcal{N}$. A Python script to compute the projection matrix P can be found in appendix A.

To enable replacement of the invalid candidate solution $\hat{\mathbf{u}}^{K,i}$ we need to define a reconstruction operator \mathcal{R} which transforms the degrees of freedom $\hat{\mathbf{p}}^{K,i}$ of the solution computed by the FVM scheme back to the ADER-DG grid. Since we have more degrees of freedom on the fine grid than on the ADER-DG grid, we can now only require that the constraints based on preservation of local averages, i.e.

$$\frac{1}{|K_\alpha|} \int_{K_\alpha} [\hat{\mathbf{u}}_h^{K,i}]_\gamma dx = \frac{1}{|K_\alpha|} \int_{K_\alpha} [\hat{\mathbf{p}}^{K,i}]_\gamma dx = [\hat{\mathbf{p}}^{K,i}]_{\alpha,\gamma} \quad (2.115)$$

for $\alpha \in \mathcal{N}_S$ and $\gamma \in \mathcal{V}$, are fulfilled in a least squares sense. We do however insist that the overall integral average

$$\int_K [\hat{\mathbf{p}}^{K,i}]_\gamma dx = \int_K [\hat{\mathbf{u}}_h^{K,i}]_\gamma dx. \quad (2.116)$$

is preserved to make sure that the scheme remains conservative.

Analogous to eq. 2.113 the left-hand side of eq. (2.115) can be simplified as follows:

$$\begin{aligned} \frac{1}{K_\alpha} \int_{K_\alpha} [\hat{\mathbf{u}}^{K,i}]_{n,\gamma} \phi_n^K dx &= \\ \int_{\hat{K}} [\hat{\mathbf{u}}^{K,i}]_{n,\gamma} \hat{\phi}_n \left(\frac{1}{N_S} \alpha + \frac{1}{N_S} \xi \right) d\xi &= \\ \sum_{n \in \mathcal{N}} \left(\prod_{d \in \mathcal{D}} ([P]_{\alpha_d, n_d}) [\hat{\mathbf{u}}^{K,i}]_{n,\gamma} \right). \end{aligned} \quad (2.117)$$

Eq. (2.116) simplifies to

$$\sum_{\alpha \in \mathcal{N}_S} \frac{1}{|K_\alpha|} [\hat{\mathbf{p}}^{K,i}]_{\alpha,v} = \frac{1}{N_S^D} \sum_{\alpha \in \mathcal{N}_S} [\hat{\mathbf{p}}^{K,i}]_{\alpha,v} = \sum_{n \in \mathcal{N}} \hat{\omega}_n [\hat{\mathbf{u}}^{K,i}]_{n,v} \quad (2.118)$$

for $v \in V$. In order to be able to carry out the computation in a dimension-by-dimension manner as for the projection, we need to solve the following constrained least squares optimization problem to find the reconstruction

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matrix $\mathbf{R} \in \mathbb{R}^{(N+1) \times N_S}$ for $\mathbf{P} \in \mathbb{R}^{N_S \times (N+1)}$ being the projection matrix and for vectors $\mathbf{p} \in \mathbb{R}^{N_S}$ and $\mathbf{u} \in \mathbb{R}^{N+1}$ representing the degrees of freedom along a coordinate axis for a single quantity:

$$\begin{aligned} & \underset{\mathbf{u} \in \mathbb{R}^{N+1}}{\text{minimize}} \quad \|\mathbf{P}_{ij} [\mathbf{u}]_j - [\mathbf{p}]_i\|^2 \\ & \text{subject to} \quad \hat{\omega}_i [\mathbf{u}]_i = \frac{1}{N_S} \sum_i [\mathbf{p}]_i. \end{aligned} \quad (2.119)$$

The corresponding Lagrange dual problem (see e.g. [14] for a general derivation) written in matrix vector notation reads

$$\begin{bmatrix} 2\mathbf{P}^T \mathbf{P} & \hat{\omega}^T \\ \hat{\omega} & 0 \end{bmatrix} \begin{bmatrix} \mathbf{u} \\ z \end{bmatrix} = \begin{bmatrix} 2\mathbf{P}^T \\ \frac{1}{N_S} \mathbf{1}^T \end{bmatrix} [\mathbf{p}], \quad (2.120)$$

where $z \in \mathbb{R}$ is a scalar Lagrange multiplier and $\mathbf{1}$ is a vector with all components equal to one. We can then compute

$$[\tilde{\mathbf{R}}] = \begin{bmatrix} 2\mathbf{P}^T \mathbf{P} & \hat{\omega}^T \\ \hat{\omega} & 0 \end{bmatrix}^{-1} \begin{bmatrix} 2\mathbf{P}^T \\ \frac{1}{N_S} \mathbf{1}^T \end{bmatrix}. \quad (2.121)$$

Since we are only interested in the first $N + 1$ rows of $\tilde{\mathbf{R}}$, the reconstruction matrix \mathbf{R} is obtained after discarding the last row, i.e.

$$[\mathbf{R}]_{0:N, 0:N_S-1} = [\tilde{\mathbf{R}}]_{0:N, 0:N_S-1}. \quad (2.122)$$

A Python notebook to compute \mathbf{R} can be found in appendix A. For more details on the transformations see [15].

Identification of Troubled Cells

Once a candidate solution in this section denoted as $\tilde{\mathbf{u}}_h^{K,i+1*}$ represented by coefficients $\hat{\mathbf{u}}^{K,i+1*}$ in a cell $K \in \mathcal{K}_h$ provided by the unlimited ADER-DG scheme is available, we can apply the set of rules defined by the troubled cell indicator to check if limiting is required. Following the terminology introduced in [3] we discriminate between two kinds of detection criteria:

1. Physical admissibility detection (PAD) can be used to incorporate domain knowledge into the simulation. Most commonly criteria of this kind are used to enforce that quantities with physical meaning remain within their respective domain of definition so that the well-posedness of the problem remains ensured. A PAD criterion is represented by a set of J inequalities of the following form:

$$\pi_j(\tilde{\mathbf{u}}_h^{K,i+1*}) > 0. \quad (2.123)$$

A candidate solution in a cell $K \in \mathcal{K}_h$ is considered admissible in a physical sense if all J inequalities are fulfilled. Considering the Euler equations as an illustrative example, if we want to ensure positivity of the pressure p and the density ρ , we will have two inequalities with $\pi_1(\tilde{\mathbf{u}}_h^{K,i+1*}) = p$ and $\pi_2(\tilde{\mathbf{u}}_h^{K,i+1*}) = \rho$.

2. Numerical admissibility detection (NAD) in the form of a relaxed version of the discrete minimum principle (DMP) ensures that the total variation of the solution remains bounded. In theory we would like to require that

$$\min_{\mathbf{y} \in V(K)} \left[\tilde{\mathbf{u}}_h^{K',i}(\mathbf{y}, t_i) \right]_v - \delta_v \leq \left[\tilde{\mathbf{u}}_h^{K,i+1*}(\mathbf{x}, t_{i+1}) \right]_v \leq \max_{\mathbf{y} \in V(K)} \left[\tilde{\mathbf{u}}_h^{K',i}(\mathbf{y}, t_i) \right]_v + \delta_v \quad (2.124)$$

holds for all $v \in \mathcal{V}$, $\mathbf{x} \in K$, $K \in \mathcal{K}_h$, i.e. the values of all quantities inside K in the candidate solution should be bounded by the solution on the Voronoi neighbors in the previous time step. The vector δ is defined as

$$\delta_v = \varepsilon \left(\max_{\mathbf{y} \in V(K)} \left[\tilde{\mathbf{u}}_h^{K',i}(\mathbf{y}, t_i) \right]_v - \min_{\mathbf{y} \in V(K)} \left[\tilde{\mathbf{u}}_h^{K',i}(\mathbf{y}, t_i) \right]_v \right) \quad (2.125)$$

for example with $\varepsilon = 10^3$ to avoid problems with floating point errors. Since it would be prohibitively expensive to compute the extrema for all polynomials, we settle for

$$\min_{\substack{K' \in V(K), \\ \alpha' \in \mathcal{N}_S}} \left[\hat{\mathbf{p}}^{K',i} \right]_{\alpha',v} - \delta_v \leq \left[\hat{\mathbf{p}}^{K,i+1*} \right]_{\alpha,v} \leq \max_{\substack{K' \in V(K), \\ \alpha' \in \mathcal{N}_S}} \left[\hat{\mathbf{p}}^{K',i} \right]_{\alpha',v} - \delta_v \quad (2.126)$$

instead, i.e. we enforce the relaxed DMP criterion only for the subcell averages on the refined equidistant grid. We apply the projection operator to obtain $\hat{\mathbf{p}}^{K,i+1*} = \mathcal{P}(\tilde{\mathbf{u}}_h^{K,i+1*})$.

The MUSCL-Hancock Scheme

For the sake of completeness we will now give a quick summary on the non-linear (in the sense of Godunov), second order accurate MUSCL³-Hancock FVM scheme. Of course this is only one possible choice for a robust, lower-order scheme that can be used as a “fallback” in the a posteriori subcell limiting approach. Following the recipe given in [2] and modified to incorporate source terms and an arbitrary number of space dimensions, the scheme consists of the following steps:

³Monotonic Upstream-Centered Scheme for Conservation Laws

1. **Compute slopes:**

$$\begin{aligned} \left[\hat{\delta}^{K,i} \right]_{d,\alpha,\gamma} = \text{minmod} \left(\left[\hat{\mathbf{p}}^{K,i} \right]_{\alpha+\mathbf{e}_d,\gamma} - \left[\hat{\mathbf{p}}^{K,i} \right]_{\alpha,\gamma}, \right. \\ \left. \left[\hat{\mathbf{p}}^{K,i} \right]_{\alpha,\gamma} - \left[\hat{\mathbf{p}}^{K,i} \right]_{\alpha-\mathbf{e}_d,\gamma} \right) \end{aligned} \quad (2.127)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \{0, 1, \dots, N_S + 1\}^D := \mathcal{N}_S^*$, $\gamma \in \mathcal{V}$, troubled grid cells $K \in \mathcal{K}_h^*$, unit vectors \mathbf{e}_d and $d \in \mathcal{D}$. A component of the index $\alpha \in \mathcal{N}_S^*$ being 0 or $N_S + 1$ implies access into $\hat{\mathbf{p}}^{K',i}$, i.e. access into the fine grid solution on the respective Voroni neighbor $K' \in V(K)$ of K . We furthermore use the common definition of the minmod function, namely

$$\text{minmod}(a, b) = \begin{cases} 0 & \text{if } ab \leq 0 \\ a & \text{if } ab > 0 \text{ and } |a| \leq |b| \\ b & \text{if } ab > 0 \text{ and } |b| < |a|. \end{cases} \quad (2.128)$$

2. **Evaluate source:**

$$\left[\hat{\mathbf{s}}^{K,i} \right]_{\alpha,\gamma} = \left[\mathbf{s} \left(\left[\hat{\mathbf{q}}^{K,i} \right]_\alpha \right) \right]_\gamma \quad (2.129)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \mathcal{N}_S^*$, troubled cells $K \in \mathcal{K}_h^*$ and $\gamma \in \mathcal{V}$.

3. **Extrapolate:**

$$\left[\mathbf{w}^{K,i} \right]_{d,e,\alpha,\gamma} = \left[\hat{\mathbf{u}}^{K,i} \right]_{\alpha,\gamma} + \frac{e}{2} \left[\hat{\delta}_d^{K,i} \right]_{\alpha,\gamma} \quad (2.130)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \mathcal{N}_S^*$, troubled cells $K \in \mathcal{K}_h^*$, $\gamma \in \mathcal{V}$, $d \in \mathcal{D}$ and $e \in \{-1, +1\} := \sigma$.

4. **Evolve:**

$$\begin{aligned} \left[\mathbf{w}^{K,i+\frac{1}{2}} \right]_{d,e,\alpha,\gamma} &= \left[\mathbf{w}^{K,i} \right]_{d,e,\alpha,\gamma} + \\ &\quad \frac{\Delta t_i}{2} \sum_{d' \in \mathcal{D}} \sum_{e' \in \sigma} \left(e' \left[\mathbf{F} \left(\left[\mathbf{w}^{K,i} \right]_{d',e',\alpha} \right) \right]_{\gamma,d'} / \left[\Delta \mathbf{x}^{K_\alpha} \right]_{d'} \right) + \\ &\quad \frac{\Delta t_i}{2} \left[\hat{\mathbf{s}}^{K,i} \right]_{\alpha,\gamma} := \\ &\quad \left[\mathbf{w}^{K,i} \right]_{d,e,\alpha,\gamma} + \frac{\Delta t_i}{2} \left[\hat{\mathbf{c}} \right]_{\alpha,\gamma} \end{aligned} \quad (2.131)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \mathcal{N}_S^*$, troubled cells $K \in \mathcal{K}_h^*$, $\gamma \in \mathcal{V}$, $d \in \mathcal{D}$ and $e \in \sigma$.

5. Solve Riemann problems:

$$\left[f^{K,i} \right]_{d,\alpha,\gamma} = \left[\mathcal{G} \left(\left[w^{K,i+\frac{1}{2}} \right]_{d,+1,\alpha-e_d}, \left[w^{K,i+\frac{1}{2}} \right]_{d,-1,\alpha+e_d}, e_d \right) \right]_{\gamma} \quad (2.132)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \mathcal{N}_{\mathcal{S}}^*$, troubled cells $K \subset \mathcal{K}_h^*$, $\gamma \in \mathcal{V}$ and $d \in \mathcal{D}$.

6. Evolve source:

$$\left[\hat{s}^{K,i+\frac{1}{2}} \right]_{\alpha,\gamma} = \left[s \left(\left[\hat{s}^{K,i} \right]_{\alpha} + \frac{1}{2} [\hat{c}]_{\alpha} \right) \right]_{\gamma} \quad (2.133)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \mathcal{N}_{\mathcal{S}}^*$, troubled cells $K \subset \mathcal{K}_h^*$ and $\gamma \in \mathcal{V}$.

7. Update solution:

$$\begin{aligned} \left[p^{L^K,i+1} \right]_{\alpha,\gamma} = & \left[p^{L^K,i} \right]_{\alpha,\gamma} - \\ & \Delta t_i \sum_{d \in \mathcal{D}} \left(\left(\left[f^{K,i} \right]_{d,\alpha+e_d,\gamma} - \left[f^{K,i} \right]_{d,\alpha,\gamma} \right) / \left[\Delta x^{K_\alpha} \right]_d \right) + \\ & \Delta t_i \left[\hat{s}^{K,i+\frac{1}{2}} \right]_{\alpha,\gamma} \end{aligned} \quad (2.134)$$

for all subgrid cells $K_\alpha \subset K$, $\alpha \in \mathcal{N}_{\mathcal{S}}^*$, troubled cells $K \subset \mathcal{K}_h^*$ and $\gamma \in \mathcal{V}$.

2.2 Profiling and Energy-aware Computing on Modern x86 Systems

2.2.1 On the Importance of Performance Profiling in Software Engineering for High Performance Computing

In a High Performance Computing (HPC) context the standard balance of design goals in Software Engineering is shifted strongly in favor of maximum application performance. In view of the ever-growing complexity of computer architectures, performance profiling has become an inevitable tool that a) provides a baseline reference on the current state of a project, b) helps to prioritize, guide and track progress of optimization efforts and c) allows comparison to other state of the art solutions as well as theoretical optima.

Modern x86 processors are equipped with one or more Performance Monitoring Units (PMU) which in conjunction with a suitable operating system provide means to obtain profiling information directly within an application. This so-called Hardware Performance Monitoring (HPM) interface allows

programming of performance event counters that in turn allow computation of metrics that are correlated with good application performance and (more recently) energy efficiency (see [16, 17, 18]). Treibig et al. illustrate best practices on how HPM can be used for performance engineering on modern multi-core processors, offer a taxonomy of common causes for suboptimal application performance and offer guidance on how profiling can be used to identify and tackle them [19].

The rest of this section is organized as follows: We first give a brief summary on the x86 Instruction Set Architecture (ISA) and its importance in current HPC systems. We then focus in more detail on features and limitations of HPM in modern x86 processors and on means how to access them in the context of a normal user-space application. The section is concluded with remarks on on-chip energy monitoring provided by some of the most recent generations of x86 processors and a literature review on findings concerning the accuracy on this feature.

2.2.2 The x86 Instruction Set Architecture and the Current Prevalence of x86 in High Performance Computing

The term x86 refers to a set of Instruction Set Architectures (ISAs) based on the philosophy of the Intel 8086 16-bit microprocessor introduced in 1978 [20]. Over the years many additions to the original ISA such as support for 32-bit (1985) and 64-bit (2003) memory addressing or the introduction of special purpose instructions e.g. for vectorized floating-point operations (MMX, SSE, AVX) or encryption (AES-NI) have been made, but nevertheless all of these changes in theory do not break backward compatibility [21]. This means that even though modern x86 processors are 64-bit capable, consist of multiple cores with complex multi-level cache hierarchies and sophisticated out of order execution pipelines, they can still correctly execute the instructions that make up a program originally written for an 8086 [20]. Due to the large number of instructions specified in a modern x86 ISA and due to the comparably complex and sometimes special purpose nature of some of them, x86 is considered to be an instance of the complex instruction set computing (CISC) paradigm (as opposed to RISC, reduced instruction set computing). Throughout this thesis we use the term “modern x86 processor” to denote a 64-bit capable multi-core processor that implements a x86 ISA that includes vector instructions for floating-point operations. Even though historically more than ten companies have produced x86 processors, nowadays this means that we refer to a processor manufactured by either Intel and AMD⁴.

⁴The market share of Intel in this “duopoly” is estimated to be between 80% and 98% in the three market segments server, desktop and laptop [22].

2. THEORETICAL CONTEXT

Architecture	Count	Percentage	Accelerator	Count	Percentage
x86	468	93.6%	None	406	81.2%
Power	23	4.6%	GPU	66	13.2%
SPARC	7	1.4%	Xeon Phi	23	4.6%
Sunway	2	0.4%	Other	5	1.0%

(a) CPU Architecture (b) Accelerator Cards

Table 2.1: Distribution of CPU architectures and accelerator cards of the supercomputers listed in the June 2016 Top500 list [23]. Systems that have both GPUs and Xeon Phi accelerator cards are listed as “Other”.

The famous TOP500 list ranks the world’s fastest supercomputers in terms of floating-point performance measured using the LINPACK benchmark. The list is published biannually in June and December by Strohmaier et al. [23]. The June 2016 edition illustrates the strong prevalence of x86 in HPC (see Table 2.1a). 468 of the 500 systems listed at the moment are based on x86 processors; 455 of them use CPUs manufactured by Intel.

A major trend directly reflected in the TOP500 list is the growing popularity of dedicated accelerator cards (see Table 2.1b). Such devices are usually connected to the main CPU via the PCI Express bus, require additional power supply and sometimes come with their own form of interconnection technology to allow for direct communication that avoids involvement of host memory or CPU. Applications that exhibit a great degree of mostly homogeneous and independent parallelism such as many types of simulations in scientific computing that in terms of runtime are dominated by vectorizable floating-point operations can greatly benefit from the presence of accelerator cards. Figure 2.1 illustrates the growing number of supercomputing systems that employ accelerators over time. Starting approximately in 2010 and 2012, respectively, two types of accelerators can be identified as the main drivers of this trend:

1. Graphics processing units used in the context of general purpose computing (GPGPU), most prominently NVIDIA devices implementing the company’s Compute Unified Device Architecture (CUDA).
2. Accelerators based on the many integrated core (MIC) paradigm, that is devices from Intel’s Xeon Phi series. These devices in essence consist of a great number (> 30) of small, interconnected x86 processors all located on a single chip and support wider vector instructions than contemporary CPUs do.

As of June 2016, 94 of the 500 machines on the list employ accelerators. Even though more recently the rate of growth in the use of accelerators seems to be slowing down, it is predicted that the demand for accelerator devices



Figure 2.1: Number of supercomputing systems on the TOP500 list [23] that employ accelerator cards between June 2006 (the time when the first such system appeared on the list) and June 2016. Systems that have both GPUs and Xeon Phi accelerators are listed as “Other”. The list is updated twice a year in June and November. The bars are centered on the day of release.

specialized in parallel floating-point computations will nevertheless remain strong [24]. This is furthermore underscored by the fact that even though only three out of the ten fastest systems today employ accelerators, nine out of the ten most energy-efficient systems on the TOP500 list (consequently leading the so-called Green500 list [25]) do.

The review of the Top500 list above clearly illustrates that at the moment x86 dominates for HPC systems on a broad scale. Considering however trends such as the use of GPUs as special purpose accelerators and the fact that the leading system “Sunway TaihuLight” is based on a custom architecture developed⁵ within China might give rise to the presumption that the market of HPC systems will become more diverse within the next couple of years. The four largest supercomputers expected to go in service in the United States by 2018 are either based on NVIDIA GPGPUs (Summit [26] and Sierra [27]) or the third generation of Intel’s Xeon Phi product line then to be used directly as a CPU (Theta and Aurora [28]).

2.2.3 Hardware Performance Monitoring in Modern x86 Processors

In 1993 Intel introduced the Pentium 60 and Pentium 66 as the earliest members of the first generation Pentium microprocessor family [29]. From doc-

⁵partly in response to export restrictions

umentation available to the general public it was well known that Pentium processors collected a lot of statistics on the interaction between running code and hardware. The means how to access the data, however, were left undocumented or required signing a nondisclosure agreement with the manufacturer. Since obtaining the metrics hinted towards in the documentation would be of great value in optimizing application performance, enthusiasts such as Terje Mathisen started to employ reverse engineering techniques and eventually managed to recover the complete set of available performance counter events [30].

Today almost all processors independent of their architecture provide some form of hardware performance monitoring (HPM) for low overhead collection of metrics describing the interaction between code and hardware [31]. The performance monitoring units (PMUs) on modern multi-core processors consist of a small number⁶ of special-purpose registers (often called module specific registers (MSRs)). There is usually one PMU dedicated to each core (more precisely hardware thread if simultaneous multithreading (SMT) is available) and additional PMUs for resources shared between cores such as the memory controller or the interconnection bus. PMU registers can be programmed freely to count the number of occurrences of hardware events such as the retirement of an instruction, a cache miss or a successful branch prediction. If used correctly they provide insight into not only where bottlenecks are in the program, but also help explain the reason why they exist [32]. The special instructions necessary to do so can only be executed with kernel privileges (ring 0) and are not part of the ISA so that for example in the x86 case there is no guaranteed backward compatibility. In Linux there are several kernel modules that allow direct or indirect access to performance counters, most prominently there are OProfile [33], perf [34] and msr. The de-facto standard library for HPM on Linux is PAPI [35], which in turn is built on top of the perf interface. Other libraries such as LIKWID [36] instead make use of msr, a module that maps the MSRs to a device file interface and due to its simplicity promises lower overheads. Documentation on the use of MSRs is available in [21, chapter 19] and [37, chapter 3.2.5] for x86 processors by Intel and AMD, respectively. In principle it would of course be possible for example in a scientific simulation code to directly use the perf or msr kernel interface to program and read performance counters, but due to in general poor documentation and the volatile nature of the syntactics and semantics of counter events across the various microprocessor architectures, generations and manufacturers this approach is usually unfeasible in practice.

In general there are two approaches towards employing HPM, namely sampling and probing [31]. The former involves periodically interrupting pro-

⁶most commonly four

program execution using system interrupts to read performance counter values and to sample the call stack which allows to correlate the obtained data with a section in the program. Overhead and inaccuracy introduced by this procedure is determined by system load and the rate at which interrupts take place. Sampling does not require modification of code or binary, but due to the unavailability of debug information and compiler optimizations such as inlining, the mapping of performance data to specific lines in code might be inaccurate. Probing, on the other hand, requires that before and after sections of interest instrumentation code is added. In essence this code programs, resets and starts the performance counters prior to the execution of the section of interest and once it has finished it reads the final counter values and stores them in a data structure ready for further analysis. The overhead of probing might be significant, especially if the part of the program to be profiled is executed very often (see [38] for a practical example). On the other hand this approach is considered to be more accurate than sampling, since it does not interrupt code execution (and thereby avoids altering the state of the processor artificially) and the procedure is only minimally invasive with respect to the generated instruction stream, i.e. the only way in which the instrumentation code can have an impact on the section of interest might be compiler optimizations across the boundaries of the section which can not be carried out any more.

We will now limit our scope to the probing approach applied in the context of modern x86 microprocessors and give an overview on overhead and accuracy of performance counters reported in literature. A great amount of research has been done over the years concerning the overhead of various ways to access performance counters in user-space applications (see for example [31, 35, 39]). Due to its great importance for the profiling infrastructure presented later in this thesis we cite Table 2.2 from [31] measuring the different overheads of reading and writing (i.e. resetting, starting and stopping) a performance counter from within the kernel, a privileged user-space application accessing the msr device file through the LIKWID library (“direct”) or an unprivileged user-space application that uses the access daemon provided by LIKWID on an Intel Haswell CPU. The latter scenario is common in the context of data centers where user applications can not run with superuser privileges but still need to use performance counters. The LIKWID daemon, for example, is available on both the SuperMUC and the CoolMUC-2 systems at Leibniz Supercomputing Center (LRZ). The overhead of libraries based on the perf interface such as PAPI is found to be slightly larger.

Literature on the accuracy of the several hundred (e.g. > 400 for Haswell) counters available in modern x86 processors, on the other hand, is limited. While metrics such as cycle count are by design accurate down to a single cycle, accuracy varies dramatically for others and error margins might be dramatic in certain scenarios. Research presented by Thomas Röhl [40] carried

Mode	Read	Write
Daemon	13144	11388
Direct	1292	656
Kernel	888	312

Table 2.2: Counter access time (median) in cycles for LIKWID on an Intel Xeon E3-1240 v3 CPU (Haswell microarchitecture). Cited from [31].

out on Intel Haswell CPUs indicates that common metrics such as branch prediction ratio, instruction retirement, memory bandwidth and load/store ratio have an average error of less or equal than 0.2%. Performance counters concerned with the multi-level caching mechanism and its complex non-uniform memory access characteristics are prone to larger error margins; an example would be L1-L2 cache transfer bandwidth with an average error $> 5\%$ in some scenarios. The undocumented counter for advanced vector extensions (AVX) floating-point instructions in general is surprisingly accurate for a range of numerical benchmarks ($< 0.05\%$), especially when compared to results from earlier microprocessor generations. Significant overcounting can happen in presence of certain AVX instructions such as the ones used to copy a subset of the values in a vector register to another.

In summary we can conclude that most common performance metrics obtained from the PMUs are highly precise and if used correctly are invaluable in HPM-assisted performance engineering. Most of the time the obtained counter values are accurate and stable enough to allow not just for qualitative but also for quantitative comparison. Since however availability, semantics and accuracy of the metrics depend to a great degree on manufacturer, generation and model of the microprocessor as well as on characteristics of the scenario under consideration (e.g. single-core vs. multi-core), a lot of factors need to be taken into consideration before valid conclusions can be drawn. The overhead introduced by layers of abstraction between counter register and user application can introduce significant overhead compared to a simple register access in certain scenarios. However due to the constant nature of the overhead and the fact that it usually does not dominate the measurement, a well-calibrated profiling instrumentation can be tuned to take overheads into account so that the adjusted metrics come remarkably close to the ones that could have been obtained from direct register access.

2.2.4 Energy Monitoring in Modern x86 Processors

The servers in common data centers and to a slightly lesser extent also in supercomputing centers almost never operate at their peak capacity all at the same time. This is the reason why today power supply to such facilities is usually underprovisioned, i.e. it is not designed with respect to the absolute

peak requirement, but rather to some worst-case average power consumption [18]. This gives rise to a more recent introduction to x86 processors: On-chip power estimation and capping capabilities.

Starting in 2011 with the Sandy Bridge microprocessor architecture Intel introduced the so-called Running Average Power Level (RAPL) interface as a way to prescribe upper bounds on the current power consumption of a CPU [18]. The following year AMD presented a similar technology called Application Power Management (APM) [41]. For the resulting embedded control task to be feasible a sophisticated on-chip model to estimate the current power consumption based on architectural events was added [42]. Both technologies expose these estimates in form of model-specific registers that are updated at a rate of about 1kHz (Intel) or 0.1kHz (AMD) [43]. In the case of Intel CPUs these MSRs contain the total amount of consumed energy as a multiple of an architecture dependent base unit (e.g. $61\mu\text{J}$ or $15.3\mu\text{J}$ for Sandy Bridge, Haswell and Haswell-EP) since the last reset of said counter. The estimates cover so-called power domains so that typically separate values for the sum over all cores, for all cores plus shared on-chip resources and for DRAM memory can be obtained. AMD CPUs, on the other hand, provide power estimates again as a multiple of an architecture defined granularity (e.g. 3.8mW for the Bulldozer microarchitecture). From now on we will use the terms “RAPL counters” to denote on-chip power estimation capabilities by both vendors.

As mentioned in the introduction energy efficiency is considered to be one of the greatest challenges in designing software for exascale machines. Reliable, low overhead on-chip estimation of CPU and DRAM energy consumption is an invaluable tool in optimizing applications towards this variable and seems to be a favorable alternative in comparison with actual measurements based on complex hardware instrumentation. Before we illustrate how RAPL counters can be used within the profiling infrastructure for ExaHyPE, we first conduct a literature review on accuracy validation, major pitfalls, subsystem modeling (from package to core) and other important observations from practice. Due to a lack of literature on AMD hardware and the fact that most of our test systems as well as those used in literature are based on these architectures, we will mostly focus on Intel Sandy Bridge and Haswell CPUs. We express the hope that accuracy bounds that hold for these architectures will act as lower bounds for future microprocessor generations.

Hackenberg et al. [43] have used external equipment including high-frequency hall effect sensors to measure power consumption at the wall outlet and the mainboard power supply of two Intel Sandy Bridge systems and one AMD Bulldozer system. For a variety of benchmarks occupying all available physical cores for eight seconds they find a small constant offset between exter-

nally measured values and those reported by the RAPL interface. The offset is negative in general, i.e. RAPL slightly underestimates the amount of consumed energy; if the DRAM power domain is included the gap closes further and correlation increases. For computationally intensive tasks the offset is slightly lower than for memory-intensive tasks. The results furthermore shows that Sandy Bridge reports an almost exact, only slightly overestimated value for an idling system; AMP on Bulldozer does not seem to compensate correctly for cores in low power states. The group highlights that measurements close to the counter update rate are challenging due to the variability in the time deltas between two consecutive updates and the lack of a precise time stamp that could compensate for it. In a more recent publication [44] Hackenberger et al. repeat their measurements for Haswell-EP and find “a significant improvement compared to RAPL on previous processor generations”. The correlation for the sum of package and DRAM RAPL is “almost perfect” and if in addition a quadratic fit is employed the deviation does not exceed 3W in any measurement. They point out that due to a new RAPL mode for the DRAM domain that is based on actual measurements the reported values are now extremely precise throughout all load scenarios.

Desrochers et al. [45, 46] instrumented a Haswell desktop system in a similar way as in [43]: Power supply is intercepted at the wall outlet, at the mainboard connector and due to their special interest in RAPL estimates for DRAM also at the respective memory banks. They again find a small constant offset in the reported RAPL CPU package values for both idle and fully loaded scenarios. If DRAM dims are idle and enter a lower power state RAPL on Haswell (unlike Haswell-EP in [43]) strongly underestimates its power consumption.

Hähnel et al. [47] use a “manually instrumented board” to compare power consumption values reported by the RAPL interface of a Sandy Bridge processor to measurements done with external equipment. They again find “that the curves’ characteristics are identical”, but measure a constant offset they attribute to the fact that they neglect DRAM power consumption altogether. They are particularly concerned with measuring short code paths ($< 5\text{ms}$). Since the RAPL counters are only updated at a rate of 1kHz with jitter of about $\pm 2\%$ this task is particularly challenging. Figure 2.2a illustrates the problem: We can not align the start of our section of interest labeled “kernel” with a counter update, in fact we do not even know precisely when the latter will happen. Similarly the point in time when the function returns will in general not be aligned with an update so that in this naive approach we would wrongly attribute energy consumption from before and after the function call to the measurement. Figure 2.2b, on the other hand, illustrates a possible solution. Before starting the actual “kernel”, we first repeatedly poll the RAPL MSR and count the number of tries that are need until a change of the counter value is observed. We then execute the kernel and



Figure 2.2: Approaches towards measuring power consumption in short code paths using RAPL counters. For the illustration we assume $t_1 - t_0$ to be of the order of milliseconds and counter updates happen at t_0 and t_1 and at unspecified times in between if $t_1 - t_0 \gtrsim 2\text{ms}$. The dark gray box indicates execution of the kernel, a light gray box labeled “p” indicates one instance of register probing. The width of the box is not true to scale and in practice we observe at the order of XXX probes before the counter value changes (compare Table 2.2). It has been shown that the approach illustrated on the right works even if the time it takes to execute the kernel is $< 1\text{ms}$. Illustration adopted from [47].

repeat the same procedure once it has finished. Via a priori calibration we can precisely determine the amount of energy that is required for one single poll simply by executing the procedure sequentially for a significant number of times and then computing the mean as the value for a single register poll. Since the number of polls we had to wait for the counter updates to happen is known we can correctly compensate for this overhead and adjust the overall energy measurement such that it only reflects the kernel function. In a case study based on single-core video decoding it was shown that if a frame is split into twelve slices and the RAPL energy consumption based on the approach explained above is measured for processing each slice individually, the sum over all slices on average differs by 1.13% compared to a measurement spanning the processing of the complete frame at once [47]. Since in ExaHyPE we will mostly be dealing with short kernel calls close to the minimum temporal resolution of RAPL energy counters we implemented this refined measurement procedure in the profiling infrastructure to be discussed in the next chapter.

With current generation x86 processors RAPL counters can not be used to obtain energy consumption estimates for individual cores. A way to get around this limitation is largely based on research tasked with predicting CPU energy consumption from a time before RAPL counters were introduced: Based on selected performance counters and other information available at runtime such as core temperature and clock speed, even relatively simple models were able to make remarkably accurate predictions. If these

models are calibrating against RAPL estimates for the complete core they can be used to predict the energy consumption of a single core.

Goel et al. [48] proposed a model based on piecewise linear regression that takes into account CPU temperature and a hardware-specific choice of the four performance counters with highest correlation to predict power for each core individually. Validation against measurements of total CPU power consumption for various x86 microprocessors introduced between 2007 and 2009 yields a median error of 1.1% - 5.2% for a diverse set of benchmark scenarios.

Also prior to the general availability of RAPL counters Takouna et al. [49] proposed a linear model based on frequency (required to be the same for all cores), squared frequency and number of active cores:

$$P(F, n) = \theta_3 F^2 + \theta_2 F + \theta_1 n + \theta_0. \quad (2.135)$$

Validation against a Nehalem-EP server with build-in CPU power measurement capabilities yields that the prediction error is less than 7% for 95% of the tested combinations of frequency F and active cores n .

Yasin et al. [50] used a particularly simple model based on normalized frequency of the individual cores of the form

$$P(c) = \theta_0 + \theta_1 \sum_{n=1}^N f_n, \quad (2.136)$$

where N is the total number of cores in the system and $f_n = F_n / F_{\max}$ is the normalized frequency of the n -th core. For a system consisting of two Sandy Bridge processors and the model calibrated against RAPL power measurements they report a mean absolute error of 4.01% and 5.23% if core power saving states (C-states, P-states) are disabled or enabled, respectively.

Even though literature in the context of RAPL counters is admittedly limited and oftentimes methodically unclear or questionable, we can still conclude that the approach of using simple linear models to predict core-level power consumption from per package aggregates is promising. In practice, however, careful calibration and extensive sanity checking is required on a case-by-case basis. Since RAPL counters on Intel platforms report total energy consumption rather than average power intake and due to the jitter in the update rate described above it is justified to assume that estimation of per-core energy consumption will be more accurate than of per-core power consumption.

Summing up the key findings of the literature review we can conclude that RAPL counters as the standard way for energy monitoring on modern x86 processors are an important tool when optimizing applications for energy

consumption. Under no circumstances do the provided estimates violate the ordering constraints of instruction streams with respect to an ordering based on real measurement with external instrumentation. They can therefore be used as a cost function to guide manual or automatic optimization. The accuracy of the counters seems to be good enough for most practical use cases; sampling frequencies at the same order of magnitude than that at which updates occur are challenging, but possible using the probing approach presented above. Further improvements with respect to temporal granularity in future architectures would nevertheless be highly beneficial. The estimation of power consumption for individual cores on a multi-core processor using simple linear models seems to be feasible and usually results in low relative errors.

Performance and Energy Profiling in ExaHyPE

In this section we will present in great detail the design and implementation of a generic performance and energy profiling infrastructure built into ExaHyPE as the main contribution covered by this thesis. The chapter is organized as follows: We will begin our discussion with an introduction to the ExaHyPE project at first focusing on high-level aspects such as where it stands in the context of current trends in Scientific Computing, what the key objectives and benefits are, which techniques and approaches it employs and what the agenda looks like with respect to past and future work. We will then give a brief summary on the architecture and usage of the ExaHyPE engine itself from a user's perspective since this defines the environment for all efforts discussed thereafter. Once this is done we can finally focus on the generic profiling infrastructure that was contributed to the project. We will again begin with some general considerations on the motivation behind the effort and what in consequence the main design goals are. After that we will illustrate the general architecture of the proposed solution, present exemplarily three implementations based on different profiling libraries and discuss ideas for future work.

3.1 The ExaHyPE Project

3.1.1 Context

Before we take a closer look at ExaHyPE, let us first consider three important aspects of contemporary scientific computing that will directly lead to the fundamental motivation that drives the project:

Simulation Pipeline: According to Bungartz et al. [51], generic simulation tasks in scientific computing can in general be subdivided into a set

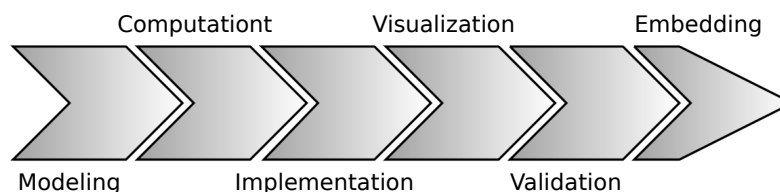


Figure 3.1: The simulation pipeline according to Bungartz et al. The process depicted is in general of iterative nature. Illustration adopted from [51].

of steps as described by the so-called simulation pipeline depicted in Figure 3.1. The pipeline model is inherently iterative in nature and describes tasks that in some form reappear in all modeling and simulation projects. For the most part all of these steps are already covered by some form of library, tool or approach with a proven track record in existing HPC applications. More often than not researchers nevertheless tend to “reinvent the wheel” and spend a significant amount of time on solving tasks that others have already worked on extensively in the past that therefore can be considered solved with respect to state of the art requirements. Most of the time side efforts such as the development of a custom linear algebra system or a data structure for adaptive mesh refinement that do not even meet the performance benchmark set by preexisting solutions. In an environment dominated by growing complexity and constrained availability of resources we would like to make the argument that the tasks as described by the simulation pipeline and subtasks thereof should be considered building blocks that need to be assembled correctly to fulfill the overall goal of such projects which is to enable new insights in the application domain. In the spirit of agile development it often makes sense to build a prototype based on existing solutions first and only after it has been determined that the overall solution does not meet the requirements to enable such insights, careful analysis and profiling needs to be applied to understand which parts are critical and need optimization. Then – only if the necessary expertise is available, of course – an effort to develop a new solution should be started and once it has been carefully validated shared with the scientific community as a building block to accelerate progress in other areas. The argument obviously has a lot of common sense built into it as this is normally the case for strong arguments, but too often in practice it is forgotten that nobody can be an expert on anything!

Exascale Computing: Simulation on supercomputers has obviously become a main driver of scientific progress. First systems capable of performing more than 10^{18} floating-point operations per second, i.e. 1 ExaFLOP/s, are expected to be operational by 2020 [52]. A subset of the

key challenges identified by Rosner et al. for the U.S. Department of Energy [53, 54] are illustrated in the following:

Energy consumption: Based on the simple extrapolation of the most energy-efficient system¹ currently listed on the Green500 list [25] an exascale machine would require about 150 MW. To put that in perspective the largest system in terms of peak power consumption² currently on the Top500 list [23] consumes less than 18 MW. Considering an average price of electricity for industrial customers in the U.S. of about 0.07 USD/kWh [55] this would result in annual expenses of more than 90M USD, a conservative estimate for Germany at 0.13 EUR/kWh would be about 170M EUR. The proposed energy target between 20 and 40 MW for an exascale system [53] gives rise to doubts expressed by many in the HPC community [56, 57] whether the assumption of such a system being introduced by 2020 is realistic. Irrespective of the exact time frame, exascale computing will induce a major shift from the perception that energy efficiency is a hardware-only challenge towards the understanding that it needs to be considered equally from both a hardware and a software perspective.

Multi-level parallelism and hybrid architectures: Since the clock frequency of chips remained more or less stagnant over the last couple of years and is even expected to decrease in an effort to reduce power consumption, increased parallelism will be the main driver of performance gains to a much stronger degree than it already is today [54]. Parallelism will therefore increase on all levels starting with instructions (pipelining, superscalar execution, VLIW) over to cores and nodes (multi-core, MIC architectures, coprocessors, accelerators) up to cross-site computation (multi datacenter computations, PRACE). Designing solutions that can make use of this trend towards “multi-level parallelism” on all levels is a great challenge both with respect to hardware and software.

Fault tolerance and resilience: If the number of cores in a chip and the number of nodes in a system increases dramatically as it will necessarily be the case in an exascale supercomputer it is an obvious observation from probability theory that the failure of components will be of significantly greater concern than in today’s petascale systems. It is therefore unavoidable to consider

¹The Shoubu supercomputer at the RIKEN research institute in Japan provides 6673.84 MegaFLOP/s per Watt.

²The Tianhe-2 system at the National Super Computer Center in Guangzhou, China, consumed 17.808 MW while running the LINPACK benchmark.

the increased probability of hardware defects in HPC software solutions that operate at this scale. Similar to common frameworks for distributed data processing and storage available today such as Google Flume [58] or Apache Hadoop [59] resilience must be a major design goal from the very beginning on.

Memory and network bandwidth bottlenecks: Data collected by Hennessy and Patterson indicates that between 1980 and 2010 processor performance increased by approximately 30% on a year by year basis. Memory bandwidth, however, only increased by about 7% annually during the same time [60]. The observation that improvements in memory technology with respect to latency and bandwidth can not keep track with advances in microprocessor performance is commonly referred to as the memory wall [61] or the memory gap [62]. A similar observation can be made for the interconnection network that provide a link between the individual compute nodes in HPC systems [63]. As a software-side answer towards these challenges induced by divergent rates of performance growth and the resulting limitations of modern hardware, novel strategies and schemes that foster even more principles such as data locality and algorithmic density must be developed to achieve scalable performance on future supercomputer architectures.

The considerations above have illustrated that there are major challenges towards exascale computing. In order to overcome them it is necessary to consider both the hardware and the software aspects of supercomputer development on equal terms. We will illustrate in chapters 3.1.2 and 3.1.3 how the ExaHyPE project directly tackles the challenges in the latter area and underscore the importance of the contributions in the context of exascale HPC.

Hyperbolic Balance Laws: The general form and a state of the art method to solve hyperbolic balance laws (HBLs) has been given in chapter 2.1. HBLs are of great interest in scientific computing since they can be used to model a variety of phenomena in physics and engineering. There are numerous areas of application including, but not limited to mechanical engineering, geophysics, electrical engineering, biology, chemistry, astrophysics, material science, civil engineering and even in economics, urban planning and the social sciences. Particularly interesting cases where numerical simulation of HBLs lead to new scientific insights reported in literature include earthquakes [64], pollutant transport [65], traffic modeling [66] and the merger of neutron stars [67]. According to Rosner et al. forefront computing problems can be divided into three categories, namely incrementally advanced computing, vo-



Figure 3.2: Logo of the ExaHyPE project [68].

racious computing and transformational computing [53]. Simulations based on HBL models can fall into all three categories. For problem scenarios that fall into the first two categories the advance to exascale computing will enable a more fine-grained understanding of the characteristics of the simulated phenomenon and allows for more accurate risk assessment at larger scales. For problems of the third category that simply can not be simulated on today's petascale machines, however, exascale systems as a result of the incorporation of novel approaches in both hardware and software will lead to new fundamental scientific findings that can really transform disciplines.

3.1.2 Vision, Objectives and Benefit

The ExaHyPE project addresses the software-side challenges posed by supercomputer development towards exascale [52]. Similar to the notion of a standardized engine commonly used in the context of computer games, the project seeks to build an engine for solving arbitrary problems formulated in terms of hyperbolic balance laws [69]. The abbreviation ExaHyPE stands for “exascale hyperbolic PDE engine”; any similarities to words in the English language are purely coincidental³.

Based on their strong expertise in areas such as numerical treatment of HBLs, large-scale simulations on some of the world's fastest HPC systems and in-depth application domain knowledge the four-year ExaHyPE project is jointly carried out by a consortium of seven institutions in Germany (Technische Universität München, Ludwig Maximilian Universität München, Frankfurt Institute for Advanced Studies, Bavarian Research Alliance GmbH), Italy (Università degli Studi di Trento), the United Kingdom (Durham University) and Russia (RSC Technologies). The consortium in addition is associated with the Leibniz Supercomputing Centre (LRZ) of the Bavarian Academy of Sciences and Humanities and cooperates with the Joint Super-

³Quotation needed...

computer Center (JSCC) of the Russian Academy of Sciences. The project was started in October 2015 and has received funding of 2.8M EUR from the European Union’s Horizon 2020 research and innovation program under grant agreement No. 671698.

The objectives of the project [68] directly reflect the challenges that are associated with the transition towards exascale computing as presented in chapter 3.1.1:

Energy efficiency on tomorrow’s supercomputer hardware: The novel algorithmic and mathematical approaches developed as part of the project are inherently designed for energy efficiency and implemented in a way such that they can easily be optimized with respect to the requirements of future energy-efficient hardware.

Scalable algorithms through dynamic adaptivity: To achieve the greatest possible accuracy in a simulation with minimal computational effort, algorithms for dynamic adaptivity will be integrated into the ExaHyPE engine by experts with decades of accumulated experience in the field and a proven track record in designing scalable solutions for adaptive refinement.

Compute-bound simulations in spite of slow memory: The novel mathematical formulations and schemes used in ExaHyPE were designed specifically with important principles such as high algorithmic density and good data locality in mind. They are particularly memory efficient and reduce the required amount of communication during parallel processing to a minimum. As a result simulations based on ExaHyPE will be able to efficiently utilize future HPC systems in spite of ever-growing network and memory performance gaps.

Extreme parallelism on unreliable hardware: Fault tolerance and resilience are key design goals of ExaHyPE. The engine will be able to dynamically schedule computations on millions of processor cores and guarantees successful completion of the simulation even if individual nodes fail in the process.

The ExaHyPE project will empower application domain experts in a multitude of fields from both academia and industry to achieve scientific progress by enabling incrementally advanced, voracious and most importantly transformational computations on future exascale systems. The engine designed by computer scientists with vast experience in the development of large-scale HPC applications combines sublime performance on a wide variety of current and future hardware architectures with ease of use so that researchers with only very little programming experience can fully exploit available resources without ever leaving their core area of expertise.

A distinctive property of the ExaHyPE project is the fact that the consortium does not only include researchers with a background in HPC and numerics, but also experts from two key areas of application, namely astrophysics and geophysics. Researchers from all contributing groups work together as equal partners in order to never lose focus on the requirements of future users of the engine. In the area of geophysics ExaHyPE will enable simulations of earthquakes and aftershocks to assess risk scenarios at unprecedented accuracy. This work will fundamentally improve our understanding of what happens during large-scale earthquakes and helps to increase “the level of preparedness [...] in expectation of the next ‘Big One’” [64]. With respect to applications in astrophysics ExaHyPE will help to tackle some of the long-standing mysteries in this field and simulations made possible by the engine are expected to enable fundamental scientific findings to help us better understand the origin of the galaxy we all live in. In 2019 ExaHyPE is to be released as an open-source package (BSD 3-Clause license) providing a user-friendly way to create state of the art, exascale ready simulation codes. This will help accelerate the pace at which application domain experts from both academia and industry can use simulations with ultimate benefits for society as a whole.

3.1.3 Approach and General Architecture

ExaHyPE is a spacetime-based engine for solving hyperbolic balance laws designed to exploit tomorrow’s exascale computing infrastructure [70, 71]. It consists of three major building blocks to be discussed in the following:

High-order space-time Discontinuous Galerkin (ADER-DG) method with a-posteriori finite volume based subcell limiting: For solving hyperbolic balance laws an ADER-DG scheme limited on a subcell level based on the MUSCL-Hancock FVM scheme as described in chapter 2.1 of this thesis is embedded into the very core of the engine. This state of the art method of arbitrary order in space and time has very favorable properties with respect to application in HPC codes: The introduction of a cell-local space-time predictor as described in chapter 2.1.7 makes the scheme inherently communication-avoiding. The fixed-point iteration that is used to obtain the latter is very favorable in terms of algorithmic density, i.e. the large ratio between the number of floating-point operations and the total memory bandwidth supports the transition from a memory-bound to a compute-bound region.

Adaptive Cartesian grids and dynamic load balancing: To minimize computational cost and to achieve the greatest possible accuracy at the same time ExaHyPE employs adaptive mesh refinement (AMR) on quadrilateral Cartesian grids. Since the engine uses the Peano framework [72] as its underlying data structure and as the main driver for

load-balanced parallelized grid iteration, technically speaking ExaHyPE itself is a “Peano application”. Peano is an open-source C++ solver framework that operates on spacetrees that employ the space-filling Peano curves for optimal cache-awareness. It can be used to solve multiple problems within one grid and has a proven track record as the foundation of numerous research projects (e.g. [73, 74, 75]) that spans over a decade.

Hardware-specific optimization of dominant compute kernels: An implementation of the ADER-DG scheme derived in chapter 2.1 essentially boils down to a number of matrix multiplications between precomputed operator matrices and matrix-valued tensor slices. The dimensionality of the appearing matrices is $(N + 1) \times (N + 1)$, where N is the polynomial degree of the ansatz functions (“the order of the scheme”), which even though arbitrarily high in theory, for practical applications reported in literature one usually has $N \in \{3, 4, \dots, 10\}$. Most libraries that provide high-performance linear algebra operations (e.g. all standard BLAS implementations) are optimized for large operands and therefore only perform unsatisfactorily for “small”⁴ matrices. ExaHyPE comes with a set of generic kernels implemented in standard C++ suitable for rapid prototyping, but for the dominating small matrix operations to achieve the highest possible performance on current and future (Intel) x86 devices, i.e. on both CPUs and accelerators, ExaHyPE can make use of the LIBXSMM library [76] to generate optimized versions of the compute kernels for the specific hardware at hand. The use of optimized kernels is essential to achieve state of the art single-core/single-thread performance. Based on the polynomial degree fixed a compile time and the types of vector instructions available on the target machine LIBXSMM directly emits assembler code which is then linked into the ExaHyPE binary. The library is developed by Intel directly, has been used successfully in major HPC projects such as SeisSol (Gordon Bell Finalist in 2014, [64]) and CP2K [77].

From a user’s perspective the workflow to create simulations based on the ExaHyPE engine is schematically illustrated in Figure 3.3. It essentially consists of the following three steps:

1. The user needs to describe the problem at hand in a simple domain specific language (DSL). Information that needs to be provided includes the size of the domain, the total time the simulation should run for, the number of variables and spatially-varying (material) parameters the problem consists of, whether shared-memory and/or distributed-

⁴Let $A \in \mathbb{R}^{M \times K}$, $B \in \mathbb{R}^{K \times N}$, $C \in \mathbb{R}^{M \times N}$ and $\alpha, \beta \in \mathbb{R}$. Then for our purposes we would for example consider the problem $C'_{mn} = \alpha A_{mk} B_{kn} + \beta C_{mn}$ small if $\sqrt[3]{MNK} \leq 80$. See github.com/hfp/libxsmm for the motivation behind this particular choice.

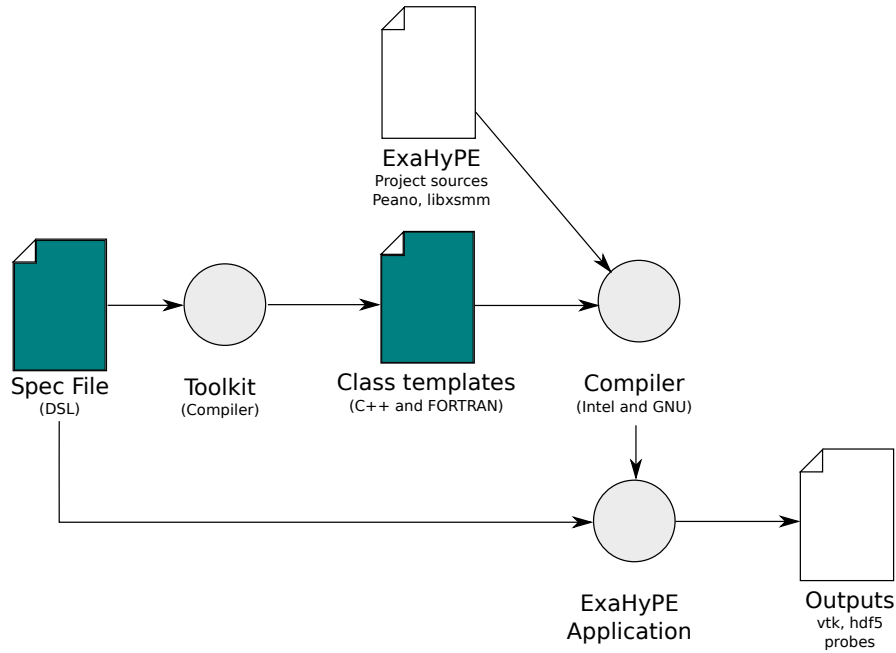


Figure 3.3: The ExaHyPE workflow as described in [78].

memory parallelization should be enabled, if and to which degree mesh refinement should take place, what the output of the simulation should look like and what language (C++ or FORTRAN) is preferred for the implementation of the problem-specific user functions. An example for a complete ExaHyPE configuration file can be found in Appendix B.

2. After the configuration file has been created it is passed to the “toolkit” application that is provided together with the actual engine. The toolkit parses the configuration file and generates glue code that is specific to both the problem and the platform under consideration. In the process stubs for all problem-specific functions such as those to prescribe eigenvalues, fluxes, sources and initial condition are created inside a folder separated from the actual ExaHyPE code.
3. The final step towards creating a complete simulation application is the implementation of problem-specific functions by the user. Only at this stage is the actual “physics” of the HBL is specified. As mentioned earlier the complete problem is described in terms of fluxes, eigenvalues, sources and initial conditions. If boundary conditions other than standard reflection or periodicity are required this can furthermore be implemented using a special API. Native interfacing with C++ and FORTRAN is supported.

Once these three steps have been completed all that is left to do is to execute the provided build script using GNU Make and to run the simulation.

A general paradigm of the ExaHyPE project is to create one specification file per experiment/machine/setup and to sacrifice some degree of flexibility in exchange for performance. This requires the user to rerun the toolkit and to recompile the application after almost all changes to the configuration, but ensures that runtime overheads are minimized and aggressive compiler optimizations can be leveraged to enable maximum performance of the resulting simulation binary. As the guidebook [69] states “ExaHyPE prefers to run the compiler more often rather than working with one or few generic executables”.

3.1.4 Current State and Next Steps

The ExaHyPE project was started in October 2015 and will span a total of four years. One year after launch, i.e. in October 2016, a prototype of the ExaHyPE engine has to be delivered that illustrates

- convergence with respect to analytic benchmarks,
- high performance of the dominant compute kernels,
- low overhead of the simulation management code, as well as
- load-balancing and scaling on a complete petascale machine.

An incomplete list of work packages for the upcoming years includes the implementation of advanced techniques with respect to fault tolerance and resilience, the integration of tooling to measure and optimize for energy consumption (based on the contributions discussed in this thesis) and provision of ways to leverage sophisticated solutions for parallel IO and visualization.

All of the results that will be presented in chapter 4 of this thesis are preliminary and given merely to discuss the implemented techniques and to illustrate their use in practice. For results on the performance of a version of the ExaHyPE engine accessible to the general public the reader must be referred to future publications.

3.2 A Generic Profiling Infrastructure for ExaHyPE

In this chapter we will discuss the main contribution of this thesis, a generic profiling infrastructure designed for extensibility and ease of use built directly into the core of the ExaHyPE engine. The section is structured as follows: First, we will begin with a quick discussion on why such tooling is necessary and what benefits it comes with. We will then focus on the key design goals as well as the resulting architecture and interface derived

thereof. Three implementations of the latter based on libraries covering various aspects of performance profiling are presented. We conclude with a quick review of ideas that might be interesting to implement in the future.

3.2.1 Motivation

As mentioned before, in essence the goal of the ExaHyPE project is to build a highly sophisticated, nevertheless easy to use engine for solving large HBL PDE problems on a variety of platforms with stellar performance. To obtain a baseline on the latter, to track progress in this area and to be able to compare it to other state of the art solutions it is essential to have sophisticated tooling for profiling available.

Besides the more traditional metrics for performance in HPC such as floating point operations per second (FLOP/s) and time to solution, in recent years it has become obvious that overall energy consumption and maximum power requirement are key metrics that large scale simulation codes need to be optimized for. The two major computing centers in Bavaria, LRZ and RRZE, are pioneers in this area whose expertise is highly valued internationally by partners from both industry and academia. The ExaHyPE team should therefore incorporate as much of their knowledge as well as experience and as many of their technologies as possible into a profiling infrastructure for ExaHyPE, but still needs to keep the tooling simple enough such that even less experienced user can gain valuable insights.

3.2.2 Design

The key design goals that were identified for a generic profiling infrastructure are listed and ways to incorporate them into a consistent design are discussed in the following:

Ease of use: The instrumentation code necessary to enable profiling for regions of interest in the application should be as simple as possible. Similar to established solutions for performance probing such as the Google Census framework [79] or LIKWID's marker API [31] we therefore chose a design that employs a tag-based system to keep track of the various regions. Independent of the actual implementation yet to be discussed, instrumentation will then look similar to the following snippet of C++ code:

```
1 MyProfilerImplementation p;  
2 p.registerTag("region1");  
3 p.start("region1");  
4 // work, work, work, ...  
5 p.stop("region1");  
6 p.writeToOstream(&std::cout);
```

Extensibility and support on a wide variety of HPC systems: In view of the plans to make ExaHyPE available as open-source software by 2019 it is essential to consider that the engine is expected to be deployed in numerous HPC systems that differ significantly not just in hardware, but also in terms of the employed software stack as well as available tools and libraries. This leads to the requirement that a profiling infrastructure for ExaHyPE needs to be designed with extensibility in mind. Researchers that require profiling to optimize their simulations in a new environment will then simply need to implement the provided profiling interface to be able to use available tools and libraries directly within ExaHyPE. Ideally this will also be a way to integrate community contributions into the project.

Low measurement overhead: The proposed generic profiling architecture for ExaHyPE forms a layer on top of existing tools and libraries. This layer is responsible to manage the assignment of individual measurements to the defined regions of interest. In order to ensure low measurement overhead it needs to be kept as thin as possible both in terms of memory and runtime requirements. Since they allow constant-time updates of stored measurement values, hash maps of carefully chosen size seem to be a suitable data structure for the task at hand.

Low maintenance effort for the ExaHyPE team: Code used for profiling is inherently specific to the hardware and software stack it was written for. In the case of hardware performance counters and model specific registers found in modern x86 processors, for example, support, naming and accuracy of events usually vary between microprocessor generations or even individual models; the consistent lag in or lack of publicly accessible documentation makes maintenance of such code a great challenge [31]. Since development of this kind of code is neither a key objective of the ExaHyPE project nor a particular area of expertise of the team, a generic profiling infrastructure must therefore be designed such that most if not all maintenance effort can be offloaded to developers of specialized libraries.

Validation and sanity checking: A profiling infrastructure for ExaHyPE should be able to make use of all available sources of information. With respect to energy profiling, for example, this means that depending on availability in a particular HPC system we want to access data from software, hardware and hybrid measurement units and whenever more than one interface is available to obtain them (e.g. perf and msr) we want to be able to switch between all of them. Only in this way can we assess the accuracy of our measurements and gain maximum confidence by employing techniques such as cross-validation and other types of sanity checking. If for a given system several ways

to obtain certain metrics are available but only one or even none has yet been implemented for use with ExaHyPE this again stresses the goal of extensibility discussed above.

Flexibility with respect to internal and external analysis: Performance metrics obtained via the generic profiling interface need to be accessible for both internal use within the engine (e.g. for optimizations at runtime) as well as external post-processing and analysis. Internally this can be ensured simply by making the object that manages profiling visible and accessible within the relevant scopes. For external analysis and post-processing the generic profiling infrastructure provides standardize export capabilities to common formats such as CSV or JSON.

3.2.3 Architecture and Interface

Figure 3.4 illustrates the general architecture of the profiling infrastructure and hints towards exemplary implementations based on different libraries for performance analysis to be discussed in the following chapter. As mentioned earlier in the discussion of design goals and ways to meet them, the virtual class `Profiler` prescribes that each concrete implementation of it needs to support a tag-based system to manage measurements with respect to predefined regions of interest in the actual simulation. In essence it describes a profiler as an object that exposes just two methods for performance measurements: `void start(const std::string& tag)` and `void stop(const std::string& tag)`. As described above, a region of interest such as a kernel call will then simply be surrounded by calls to the two methods. For the sake of decreased overhead during the actual measurement, before a tag can be used for profiling, depending on the actual implementation, it either should or has to be registered via `void registerTag(const std::string& tag)`. Registration of a tag in the various implementations that are presented exemplarily in the following allocates the necessary field in one or more hash maps (e.g. `std::unordered_map<std::string, value-type>` in C++) and initializes the fields to zero for aggregated values. In general the interface does not prescribe if a profiler has to report aggregated or individual measurement values. Registering tags in advance reduces the one-time overhead when the first measurement for this tag is started.

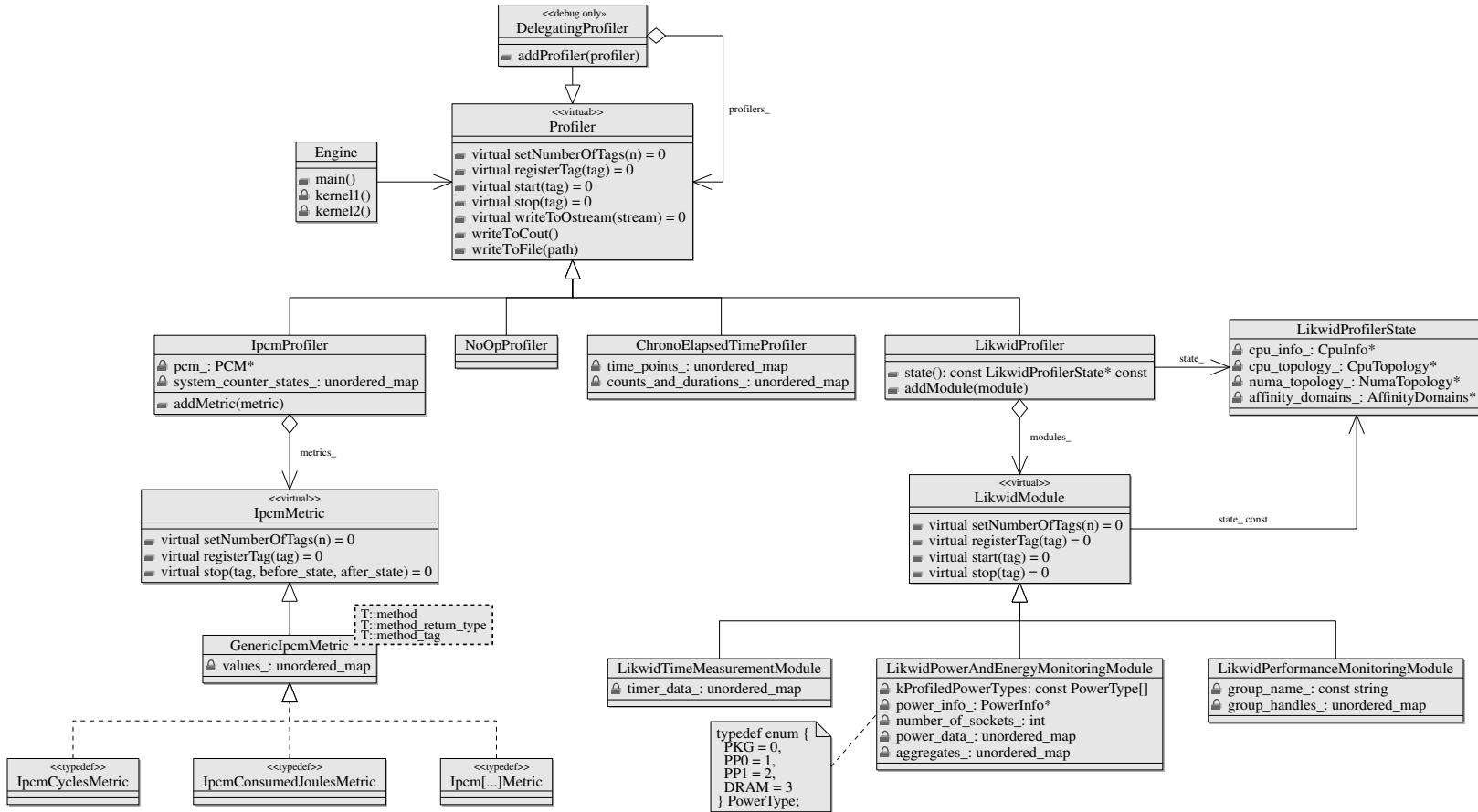


Figure 3.4: Class diagram illustrating the proposed generic profiling architecture and exemplary implementations based on different libraries for performance analysis.

Integration of the profiling infrastructure into the overall project has been carried out as follows: A new optional block in ExaHyPE's domain specific configuration language is used to enable profiling and to specify the desired profiler type and metrics (see appendix B for details). Profiling is disabled by default. Furthermore, a member of type `std::unique_ptr<Profiler>` was added to the engine's `Solver` class. During initialization of the solver and based on the parsed configuration file available at runtime a `ProfilerFactory` creates the desired type of profiler. The solver object takes ownership of the profiler, i.e. the existence of the profiler is tied to the life time of the solver. It is responsible to register the required tags, e.g. one for each kernel in the solver. If profiling is enabled, the toolkit is used to add the necessary instrumentation that calls the profiler to start and stop the measurements. It simply becomes part of the glue code that needs to be generated either way to connect the ExaHyPE kernels and the problem-specific functions provided by the user.

After the simulation has finished the profiler can export the collected data via a call to `void writeToOstream(std::ostream* stream)` to an arbitrary output stream. JSON (JavaScript object notation) has been chosen as the default output format due to its simplicity, flexibility and widespread support by almost all common tools that might be used for analysis and visualization of the data. Export to the even simpler CSV (comma-separated values) format is available as well. If required in the future it would be trivial to implement export to other file formats such as XML (extensible markup language).

3.2.4 Exemplary Implementations

As part of this thesis three implementations of the profiling interface discussed in the previous section were developed and are to be presented in the following.

STL Chrono

A relatively trivial implementation of the interface called `ChronoElapsed-TimeProfiler` makes use of the `chrono` library that is part of the C++ Standard Template Library (STL). It can be used to count the number of times a particular region in the code has been executed and to measure the overall wall clock time spent therein. Since an implementation of the STL is shipped with all operating systems and C++ compilers, the profiler does not have any external dependencies. The accuracy of `chrono` depends on the specific combination of compiler, operating system and hardware; it is usually at the order of nanoseconds. The overhead induced by profiler and library is analyzed at startup; results are presented in chapter 4.

Intel Performance Counter Monitor

The Intel Performance Counter Monitor (PCM) library [80] provides access to RAPL and performance event counters⁵ from within a user-space application in an easy to use way. The library only supports Intel CPUs and is available as open-source software under the same BSD3-Clause license as ExaHyPE. The class `IpcmProfiler` that is shipped as part of ExaHyPE is a thin layer on top of PCM that is responsible for initialization, deinitialization and management of the mapping between individual measurements and regions of interest. An instance of `IpcmProfiler` owns a set of `IpcmMetric` objects as specified by the user in the configuration file. Each implementation of `IpcmMetric` can be mapped to exactly one of the metrics that the library provides. In PCM a metric is represented by a function of the form `double/int get...(SystemCounterState& before, SystemCounterState& after)`, where `...` is a placeholder for the name of the actual metric. Due to this particularly homogeneous structure all provided `IpcmMetric` implementations are simply instantiations of a single C++ template called `GenericIpcmMetric`. See Table 3.1 for the subset of metrics provided by version 2.11 of the Intel PCM library that are accessible via `IpcmProfiler` in ExaHyPE.

A binary that uses the PCM library to collect performance metrics is required to be executed with superuser privileges or needs the accompanying `pcm-sensor` daemon to be accessible. In consequence PCM-based profiling does not work on LRZ's SuperMUC and CoolMUC-2 by default at the moment. A minor drawback of using PCM in conjunction with ExaHyPE is the fact that the library uses exceptions, but exceptions are disallowed in ExaHyPE. The author has created a patched version of PCM 2.11⁷ so that the library does not throw exceptions but aborts the application instead⁸. At startup `IpcmProfiler` performs a set of measurements to estimate the overhead introduced by profiling.

⁵PCM uses the `msr` kernel module on Linux and a custom driver on Windows.

⁶Modern x86 processors employ speculative execution, i.e. they execute more instructions than absolutely necessary in an effort to accelerate the overall computation. An illustrative example might be a situation where the CPU encounters a conditional jump instruction. If the condition depends on a value which must be fetched from memory and in consequence would stall execution, the CPU may decide to execute both possible branches in the meantime and once the condition can be evaluated only the instructions on the correct branch are "retired".

⁷The patched version is available at git.io/vis7G.

⁸This really does not change anything in our case, since the library only throws exceptions if performance analysis does not work (e.g. due to lack of superuser privileges) and in these scenarios there really is not much we can do to recover from the exception either way.

3.2. A Generic Profiling Infrastructure for ExaHyPE

Value Type	PCM Function	Explanation
double	getCycles	Number of core clock cycles
uint64	getInstructionsRetired	Number of retired instructions ⁶
double	getIPC	Average number of retired instructions per cycle
uint64	getBytesReadFromMC	Number of bytes read from DRAM memory controllers
uint64	getBytesWrittenToMC	Number of bytes written to DRAM memory controllers
uint64	getL2CacheHits	Number of L2 cache hits
uint64	getL2CacheMisses	Number of L2 cache misses
uint64	getL3CacheHits	Number of L3 cache hits
uint64	getL3CacheMisses	Number of L3 cache misses
double	getCyclesLostDue-L2CacheMisses	Estimate of how many cycles were lost due to L2 cache misses
double	getCyclesLostDue-L3CacheMisses	Estimate of how many cycles were lost due to L3 cache misses
double	getConsumedJoules	Energy consumed by processor (excl. DRAM)
double	getDRAMConsumedJoules	Energy consumed by DRAM

Table 3.1: Metrics provided by version 2.11 of the Intel PCM library [80] accessible via `IpcmProfiler` in ExaHyPE.

LIKWID

Without doubt the most interesting implementation of the `Profiler` interface is based on LIKWID. While the authors originally described LIKWID, short for “Like I Knew What I’m Doing”, as “a set of performance-related command line tools targeting X86 processors” [31], today it can be considered a full-fledged library for application profiling. Even though the C-API available since version 4.0 feels hard to use from a subjective point of view⁹ and examples are still scarce, it exposes all of LIKWID’s features related to areas such as performance counters, topology analysis and execution control (e.g. “thread pinning”). The framework is under active development and maintained by researchers at the Erlangen Regional Computing Center (RRZE). It is designed to be as lightweight as possible and optimized for low overheads.

The class `LikwidProfiler` acts as a thin adapter between LIKWID and the ExaHyPE engine. The profiler consists of one or more modules that all implement the `LikwidModule` interface. A module directly mirrors all of the

⁹Within the scope of this thesis changes by the author have been accepted into version 4.1.2 of LIKWID to improve the usability of the C-API in a C++ application.

functionality of the respective Likwid C-API module. At the moment the API consists of the following five modules:

Time measurement module: This module exposes low-overhead functionality to measure how long it takes to execute a certain region of interest both with respect to wall clock time or the number of CPU cycles.

Power and Energy monitoring module: Based on the msr kernel module this part of the API provides convenient access to the processor's RAPL counters to obtain estimates for power and energy consumption of the CPU.

Performance monitoring module: This module allows programming and readout of the various PMUs that are part of modern x86 CPUs. Similar to the event sets defined in PAPI, the developers of LIKWID provide a set of predefined groups which are designed to be independent of underlying hardware-dependent events with respect to details such as exact naming, semantics and availability. Concerning RAPL counters the API is somewhat redundant in the sense that both the performance monitoring module as well as the power and energy monitoring module offer the same (but in case of the latter more convenient to use) functionality.

Thread affinity module: Servers in an HPC context are usually more complex than normal desktop computers with respect to NUMA characteristics as well as number and kind of available compute resources (multi-core CPUs, accelerators, etc.). The thread affinity module exposes functionality to simplify management of this degree of complexity, to allow for convenient logical addressing of the available resources and to restrict execution of an application to certain parts of the hierarchy (e.g. pinning to a specific hardware thread).

Thermal monitoring module: This module provides access to MSRs holding static (e.g. thermal design point, maximum core temperature) and dynamic (e.g. current temperature in a core) values relevant to thermal management of the CPU. Temperature sensors had been available in microprocessors long before any other form of hardware monitoring was introduced. Since temperature can be seen as an indirect measure for running average power consumption with very wide sampling width, historically it has played an important role in the modeling of CPU power consumption.

The `LikwidProfiler` class is a very thin layer on top of Likwid that manages initialization and finalization of the various Likwid modules, hides some of the rough edges of the relatively new C-API in a reasonably robust piece of C++ code and manages the export of the collected data. The class `LikwidPerformanceMonitoringModule` makes use of the predefined event

sets to collect performance metrics in a platform-agnostic way. Amongst others, at the moment the following groups are available within ExaHyPE: FLOPS_AVX, BRANCH, CLOCK, DATA, ENERGY, ICACHE, L2, L2CACHE, L3, L3CACHE, MEM. For the sake of brevity we omit a detailed discussion of the provided metrics and refer to the respective online documentation¹⁰ instead. The library needs either root privileges or its own daemon running with root privileges to be able to collect performance metrics. The corresponding daemon `likwid-accessD` is available on SuperMUC and CoolMUC-2. This type of indirect access to performance data induces overhead as a result of the employed socket file communication mechanism and additional security checks by the daemon necessary to filter out dangerous or malicious requests. The profiler implementation will give estimates of the latter at startup so that data can be recalibrated accordingly. Likwid supports modern x86 CPUs by Intel and AMD as well as Intel Xeon Phi accelerators.

3.2.5 Future Work

An interesting area for future work will be the extension of the generic profiling infrastructure to parallel ExaHyPE runs both with respect to shared-memory and distributed-memory parallelization. In this context, however, most work will need to be done within Peano rather than ExaHyPE. Careful evaluation regarding preexisting and overlapping functionality as well as possible synergies must proceed all efforts in this direction. The end goal should be a profiling infrastructure which enables insights across all levels of the distributed application so that a complete picture can be used to guide optimization efforts targeted at performance and energy efficiency of the whole simulation.

At the moment the available profiling infrastructure is used exclusively for a posteriori analysis. A promising idea for future work would be to use the available metrics directly within the engine, e.g. for tasks such as load-balancing based on energy consumption.

In an effort to increase confidence in the available metrics and to increase the number of HPC systems for which the engine can offer profiling out of the box, it will be worthwhile to implement support for additional performance measurement libraries. Intel PCM and LIKWID both use the same `msr` kernel module to obtain data; it would for example be interesting to include libraries based on `perf` and `PAPI`. Nowadays almost all supercomputing centers offer ways to obtain real hardware measurements for energy-related metrics. Support for such interfaces would be of great value in the assessment of on-chip metrics. The proposed interface could furthermore be used to add instrumentation for libraries that go beyond classical CPU performance profiling such as Valgrind [81] or MeterPU [82].

¹⁰<https://github.com/RRZE-HPC/likwid/tree/master/groups>

Chapter 4

Preliminary profiling results, case studies

Overhead: measure wall clock time of sleep(0, 1, 2, ...)

System: Coolmuc, HW, normal frequency, turbo disabled, HT disabled?

Best practices: Performance patterns

- Load imbalance
- **Bandwidth saturation** (main memory, L3)
- **Strided or erratic data access** (“Cache-based architectures require contiguous data accesses to make efficient use of bandwidth due to the cache line concept”. badly ordered loop nests, inappropriate data structure)
- **Bad instruction mix** (compiler, degree of vectorization, expensive operations (sqrt, divide))
- **Limited instruction throughput** (load, multiply), related to previous
- **Microarchitectural anomalies:** Alignment, ...
- Synchronization overhead
- False cache line sharing
- **Bad page placement on ccNUMA** (pinning)

HPM groups:

- Memory bandwidth close to peak
- Low bandwidth, high store/load counts, program memory bound, cache ratios
- Ratio FP / instructions retired

4. PRELIMINARY PROFILING RESULTS, CASE STUDIES

- Static analysis: High pressure on single unit
- Hardware specific alignment things
- System vs. user time?!
- Analytic benchmark: Introduction, derivation
- Pie-chart per kernel
- Case-studies HPM-based optimization efforts: Cache-misses, compile-time (\rightarrow Toolkit philosophy)
- Degree \rightarrow Wallclock, Energy (AMR)
- Static mesh $\Delta x \rightarrow$ Error for polynomials (convergence tables)

Chapter 5

Conclusion and Outlook

- PA is important
- ExaHyPE as an answer to exascale challenges
- Applications

Chapter 6

Acknowledgment

Appendix A

Computation of the Discrete ADER-DG Operators

A. COMPUTATION OF THE DISCRETE ADER-DG OPERATORS

```
In [1]: import numpy as np
np.set_printoptions(precision=3);
import sympy as sp
sp.init_printing(use_latex=True);
x = sp.symbols("x");

## Settings
N = 4; # Degree of the ADER-DG scheme in space and time
N_S = 2*N + 1; # Number of subcell averages on the fine grid
##
```

Legendre Polynomials

$$P_0(x) = 1, P_1(x) = x, P_{n+1}(x) = \frac{1}{n+1}[(2n+1)xP_n(x) - nP_{n-1}(x)]$$

```
In [2]: def NextLegendrePolynomial(n, P_, P__):
P = ((2*n + 1) * x * P_ - n * P__) / (n+1);
return P.simplify();

P = []; P.append(1); P.append(x);

for i in range(1, N+1):
P.append(NextLegendrePolynomial(i, P[i], P[i-1]));

P = P[N+1];
P
```

```
Out[2]:  $\frac{x}{8}(63x^4 - 70x^2 + 15)$ 
```

Gauss-Legendre nodes $\tilde{\xi}$

```
In [3]: xi_tilde_sym = sorted(sp.solve(P, x));
xi_tilde_sym
```

```
Out[3]:  $\left[ -\sqrt{\frac{2\sqrt{70}}{63} + \frac{5}{9}}, -\sqrt{-\frac{2\sqrt{70}}{63} + \frac{5}{9}}, 0, \sqrt{-\frac{2\sqrt{70}}{63} + \frac{5}{9}}, \sqrt{\frac{2\sqrt{70}}{63} + \frac{5}{9}} \right]$ 
```

Gauss-Legendre nodes $\hat{\xi}$ on $[0, 1]$

```
In [4]: xi_hat_sym = [((xi+1) / 2).simplify() for xi in xi_tilde_sym];
xi_hat = np.array([xi.evalf() for xi in xi_hat_sym]).astype(np.float64);
print xi_hat;

[ 0.047  0.231  0.5    0.769  0.953]
```

Gauss-Legendre weights $\hat{\omega}$ on $[0, 1]$

```
In [5]: A = np.matrix([xi_hat**i for i in range(0, N+1)]);
b = np.array([sp.integrate(x**i, (x, (0, 1))).evalf()
              for i in range(0, N+1)]).astype(np.float64);

omega_hat = np.linalg.solve(A, b);
print omega_hat;

[ 0.118  0.239  0.284  0.239  0.118]
```

Lagrange interpolation polynomials L_i on $[0, 1]$ with nodes $\hat{\xi}$

$$L_i(\xi) = \prod_{j \neq i} \frac{\xi - \hat{\xi}_j}{\hat{\xi}_i - \hat{\xi}_j}, i = 0, \dots, N$$

```
In [6]: def L(i, xi_hat):
        xi_skip = xi_hat[:i] + xi_hat[i+1:];
        numerator = sp.prod([x - xi for xi in xi_skip])
        denominator = sp.prod([xi_hat[i] - xi for xi in xi_skip])
        return numerator / denominator

psi = [sp.lambdify(x, L(i, xi_hat_sym)) for i in range(N+1)]

In [7]: def dL(i, xi_hat):
        return sp.diff(L(i, xi_hat))

dpsi = [sp.lambdify(x, dL(i, xi_hat_sym)) for i in range(N+1)]
```

Left Reference Element Flux Operator l

```
In [8]: l = np.array([psi[i](0.0) for i in range(0, N+1)]);
print l

[ 1.551 -0.893  0.533 -0.268  0.076]
```

Right Reference Element Flux Operator r

```
In [9]: r = np.array([psi[i](1.0) for i in range(0, N+1)]);
print r

[ 0.076 -0.268  0.533 -0.893  1.551]
```

Right Reference Element Mass Operator R

```
In [10]: R = np.fromfunction(np.vectorize(
    lambda i, j: psi[i](1.0) * psi[j](1.0)), (N+1, N+1), dtype=int);
print R
```

```
[[ 0.006 -0.02  0.041 -0.068  0.118]
 [-0.02  0.072 -0.143  0.239 -0.416]
 [ 0.041 -0.143  0.284 -0.476  0.827]
 [-0.068  0.239 -0.476  0.798 -1.386]
 [ 0.118 -0.416  0.827 -1.386  2.407]]
```

Reference Element Stiffness Operator K

```
In [11]: K = np.fromfunction(np.vectorize(
    lambda i, j: omega_hat[j] * dpsi[i](xi_hat[j])), (N+1, N+1),
    dtype=int);
print K
```

```
[[ -1.201 -0.46  0.171 -0.117  0.131]
 [ 1.825 -0.363 -0.817  0.444 -0.464]
 [-0.958  1.15  0.    -1.15  0.958]
 [ 0.464 -0.444  0.817  0.363 -1.825]
 [-0.131  0.117 -0.171  0.46  1.201]]
```

Iteration Matrix \tilde{K}

```
In [12]: Ktilde = np.linalg.inv(R - K);
print Ktilde
```

```
[[ 0.53 -0.124  0.086 -0.066  0.044]
 [ 1.039  0.559 -0.151  0.102 -0.066]
 [ 1.007  1.022  0.57 -0.151  0.086]
 [ 0.981  1.016  1.022  0.559 -0.124]
 [ 1.017  0.981  1.007  1.039  0.53 ]]
```

Projection Operator P

```
In [13]: P = np.fromfunction(np.vectorize(
    lambda i, j:
        sum([omega_hat[k] *
            psi[j](1.0/N_S * i + 1.0/N_S * xi_hat[k])
            for k in range(0, N+1)])), (N_S, N+1), dtype=int);
print P
```

[9.472e-01	5.620e-02	-2.341e-03	-1.894e-03	8.191e-04]
[2.055e-01	9.434e-01	-2.174e-01	9.415e-02	-2.558e-02]
[-5.363e-02	8.698e-01	2.435e-01	-7.997e-02	2.030e-02]
[-6.623e-02	4.256e-01	7.658e-01	-1.644e-01	3.913e-02]
[-1.333e-03	1.087e-02	9.809e-01	1.087e-02	-1.333e-03]
[3.913e-02	-1.644e-01	7.658e-01	4.256e-01	-6.623e-02]
[2.030e-02	-7.997e-02	2.435e-01	8.698e-01	-5.363e-02]
[-2.558e-02	9.415e-02	-2.174e-01	9.434e-01	2.055e-01]
[8.191e-04	-1.894e-03	-2.341e-03	5.620e-02	9.472e-01]

Reconstruction Operator R

```
In [14]: m1 = np.append(2*P.T.dot(P), [omega_hat], axis=0);
m1 = np.append(m1, np.append(omega_hat, 0.0).reshape(N+2, 1), axis = 1)

m2 = np.append(2*P.T, [1.0/N_S * np.ones(N_S)], axis=0)

Rtilde = np.linalg.solve(m1, m2);
R = np.delete(Rtilde, N+1, 0)
print R
```

[1.014	0.113	-0.13	-0.073	0.04	0.074	0.007	-0.075	0.03]
[-0.064	0.514	0.468	0.195	-0.041	-0.112	-0.021	0.095	-0.034]
[0.038	-0.137	0.066	0.32	0.426	0.32	0.066	-0.137	0.038]
[-0.034	0.095	-0.021	-0.112	-0.041	0.195	0.468	0.514	-0.064]
[0.03	-0.075	0.007	0.074	0.04	-0.073	-0.13	0.113	1.014]

Appendix B

A Sample ExaHyPE Configuration File

```
1  /**
2   * This file is part of the ExaHyPE project.
3   * Copyright (c) 2016 http://exahype.eu
4   * All rights reserved.
5   *
6   * The project has received funding from the European
7   * Union's Horizon 2020 research and innovation
8   * programme under grant agreement No 671698. For
9   * copyrights and licensing, please consult the
10  * webpage.
11  *
12  * Released under the BSD 3 Open Source License.
13  * For the full license text, see LICENSE.txt
14  */
15
16 /**
17  * Euler Flow. A simple project.
18  */
19
20 exahype-project Euler
21   peano-kernel-path = ./Peano
22   exahype-path      = ./ExaHyPE
23   output-directory  = ./ApplicationExamples/EulerFlow
24   architecture      = noarch
25
26   computational-domain
27     dimension = 2
28     width     = 1.0, 1.0
29     offset    = 0.0, 0.0
30     end-time  = 0.12
31   end computational-domain
32
33   shared-memory
34     identifier    = autotuning
```

B. A SAMPLE ExaHyPE CONFIGURATION FILE

```
35     cores                = 4
36     properties-file = sharedmemory.properties
37 end shared-memory
38
39 distributed-memory
40     identifier = static_load_balancing
41     configure  = {hotspot,fair,ranks_per_node:4}
42     buffer-size = 64
43     timeout     = 120
44 end distributed-memory
45
46 optimisation
47     fuse-algorithmic-steps      = on
48     fuse-algorithmic-steps-factor = 0.99
49 end optimisation
50
51 profiling
52     profiler = LikwidProfiler
53     metrics  = {LikwidPowerAndEnergyMonitoringModule}
54     likwid_inc = /usr/local/include
55     likwid_lib = /usr/local/lib
56 end profiling
57
58 solver ADER-DG MyEulerSolver
59     variables      = 5
60     parameters     = 0
61     order          = 3
62     maximum-mesh-size = 0.15
63     time-stepping  = global
64     kernel         = generic::fluxes::nonlinear
65     language       = C
66
67     plot vtk::Cartesian::cells::ascii
68         variables = 5
69         time      = 0.0
70         repeat    = 0.05
71         output    = ./conserved
72         select    = {}
73     end plot
74 end solver
75 end exahype-project
```

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