

Athena X-IFU synthetic observations of galaxy clusters to probe the chemical enrichment of the Universe

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ABSTRACT

Answers to the metal production of the Universe can be found in galaxy clusters, notably within their intra-cluster medium (ICM). The X-ray Integral Field Unit (X-IFU) on board the next-generation European X-ray observatory *Athena* (2030s) will provide the necessary leap forward in spatially-resolved spectroscopy required to disentangle the intricate mechanisms responsible for this chemical enrichment. In this paper, we investigate the future capabilities of the X-IFU in probing the hot gas within galaxy clusters. From a test sample of four clusters extracted from cosmological hydrodynamical simulations, we present comprehensive synthetic observations of these clusters at different redshifts (up to $z \leq 2$) and within the scaled radius R_{500} performed using the instrument simulator SIXTE. Through 100 ks exposures, we demonstrate that the X-IFU will provide spatially resolved mapping of the ICM physical properties with little to no biases ($\lesssim 5\%$) and well within statistical uncertainties. The detailed study of abundance profiles and abundance ratios within R_{500} also highlights the power of the X-IFU in providing constraints on the various enrichment models. From synthetic observations out to $z = 2$, we have also quantified its ability to track the chemical elements across cosmic time with excellent accuracy, and thereby to investigate the evolution of metal production mechanisms as well as the link to the stellar initial mass-function. Our study demonstrates the unprecedented capabilities of the X-IFU of unveiling the properties of the ICM but also stresses the data analysis challenges faced by future high-resolution X-ray missions such as *Athena*.

Key words. galaxies: clusters: intracluster medium – galaxies: abundances – galaxies: fundamental parameters – techniques: imaging spectroscopy – methods: numerical – X-rays: galaxies: clusters

1. Introduction

Metals and other heavy elements in the intra-cluster medium (ICM) represent a fossil record of the chemical evolution of the Universe. Trapped in the dark matter (DM) potential of galaxy clusters (White et al. 1993), they remain unaltered within the optically-thin, collisionless thermal plasma. Elements originate within stars or through supernovae (SN), before being spread by stellar winds or by the SN explosions. Hence, the chemical enrichment of a given cluster relates to the integrated star formation history of the cluster, as well as to the overall stellar initial mass function (IMF). The abundances and spatial distribution of metals in the ICM can also be connected to its dynamical history and to the mechanical action of active galactic nuclei (AGNs) outflows or jets (e.g. Gaspari et al. 2011).

Most of the low-mass elements (C, O, Mg, Si, and S) are produced by end-of-life massive stars ($\geq 10 M_{\odot}$) undergoing core-collapse supernovae (SN_{cc}; see Nomoto et al. 2013, for a review). The evolution of SN_{cc}-related enrichment through time is dictated by the initial mass and metallicity of the progenitor star. High-mass elements, from Si-like elements (Al, Si, S, Ca, and Ar) to Fe and Ni, are on the other hand the result of thermonuclear reactions occurring during the explosion of white dwarfs (type Ia supernovae – SN_{Ia}; Hillebrandt et al. 2013). Although the mechanisms of these explosions – either via accretion of a companion star onto the white dwarf (Whelan & Iben 1973) or via mergers of binary systems (Webbink 1984) – is still poorly understood (see Maoz et al. 2014), the timescale of these events, related to longer-living low-mass stars, suggests a later enrichment across cosmic time. Traces of other elements (C, N, Ne, and Na) can also be produced when low- and

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intermediate-mass stars (typically $\leq 6 M_{\odot}$) enter their asymptotic giant branch (AGB) phase (Iben & Renzini 1983). The individual study of these phenomena based on detailed observations of nearby SN is difficult as they are very rare. Rather than a direct study on stellar populations, the detailed spectroscopic study of the ICM is an interesting alternative probe to test metal production models up to the early periods of the Universe.

Beyond the first steps in high-resolution X-ray spectroscopy (Canizares et al. 1979, 1982) and despite the lack of spatial resolution (Peterson & Fabian 2006), the advent of high-resolution grating instruments such as *XMM-Newton*/RGS (den Herder et al. 2001) and *Chandra*/HETG (Canizares et al. 2005) drastically changed our view of the ICM enrichment, by giving access for the very first time to a large number of atomic lines (de Plaa et al. 2007; de Plaa 2013; Molendi et al. 2016; Werner et al. 2007). Clusters have, therefore, become excellent laboratories in which to test plasma physics and the chemical enrichment models up to the present epoch (see Werner et al. 2008, for a review). Despite limited spectral resolutions, instruments based on charged coupled devices (CCDs) have also been pushed to the maximum of their abilities to benefit of their spatial resolution in investigating the spatial distribution of chemical elements in the ICM (de Grandi & Molendi 2009; Mernier et al. 2016a,b, 2017).

The perspective of micro-calorimeter-based imaging spectrometers, such as the soft X-ray spectrometer (SXS) on board *Hitomi* (Takahashi et al. 2016), has opened new possibilities in studying the ICM: from the spatial scales of the enrichment (sources of production, processes of mixing and dispersion) to the kinematics of the hot gas (turbulence, shocks Hitomi Collaboration 2016, 2018a,b,c), which complement the indirect estimates via surface brightness and warm gas tracers (e.g. Churazov et al. 2012; Gaspari & Churazov 2013; Hofmann et al. 2016; Gaspari et al. 2018). Unfortunately, the short lifetime of the SXS gave only a glimpse of its potential. These renewed capabilities in galaxy-cluster observation now rely on future missions, such as the X-ray Recovery Imaging and Spectroscopy Mission (XRISM; Ishisaki et al. 2018) or the Advanced Telescope for High-ENergy Astrophysics (*Athena*; Nandra et al. 2013). Namely, the X-ray Integral Field Unit (X-IFU) on board the future European X-ray observatory (Barret et al. 2016; Pajot et al. 2018), will provide narrow-field observations ($5'$ in equivalent field-of-view diameter) over the 0.2–12 keV bandpass, with a required $5''$ spatial resolution and an unprecedented spectral resolution of 2.5 eV (required up to 7 keV).

Investigating the chemical enrichment of the Universe is one of *Athena*'s prime science objectives (Ettori et al. 2013; Pointecouteau et al. 2013) which drives top-level performances of the telescope. In addition to the spectral resolution of the X-IFU, which will allow to resolve faint atomic lines of less abundant elements, this science objective drives the need for a high effective area of the telescope along with a well-calibrated low energy band, required to accurately resolve lines of light elements such as C (≥ 0.2 keV). Number of breakthroughs on the study of chemical species and their evolution should in fact come from measurement in the low-energy band, where the effective area is the highest. The fine spectroscopic capabilities of the X-IFU in this energy band will probe the production and circulation of metals within galaxy clusters across cosmic time, up to a redshift of $z \leq 2$ and a distance of R_{500}^1 from the cluster's centre. By accurately measuring the

abundances of the most common elements (e.g. O, Si, S, and Fe), the X-IFU will be capable of constraining the number of time-integrated SN_{Ia} and SN_{cc} products. For the first time, the spatially-resolved measurements of less abundant elements (e.g. C, Al, S, and Ca) as well as rare elements (e.g. Mn, Cr, and Ti) will provide insights on the initial metallicity of the SN_{Ia} progenitors, and therefore on their formation mechanisms. The science of the chemical enrichment is a driver of the performance of the instrument, which needs to be assessed before launch.

In this paper, we investigate the feasibility of recovering the physical parameters of the ICM through X-IFU observations. Careful attention is given to the different enrichment mechanisms and their evolution over time. We used a sample of four simulated galaxy clusters with different masses studied at different redshifts, obtained via hydrodynamical cosmological simulations (Rasia et al. 2015; Biffi et al. 2017). These objects are passed as input to a dedicated end-to-end (E2E) simulation pipeline of the X-IFU instrument, based on the simulator SIXTE (Wilms et al. 2014). In Sect. 2, we present the properties of the sample of simulated clusters. This is followed by a detailed description of our simulation pipeline (Sect. 3). The data analysis, post-processing procedures and results validation are in turn described (Sect. 4). The outputs of our synthetic observations obtained through the pipeline for the four local clusters are then used (Sect. 5) to infer the main properties of the sample and study its enrichment. This investigation is also extended to higher redshift values (Sect. 6) to look into the X-IFU abilities to capture the evolution of abundances through cosmic time. Finally, results and outcomes of our study are discussed (Sect. 7).

2. Generation of the cluster sample

The sample of four clusters of galaxies analysed in this study is taken from Biffi et al. (2018) and includes two massive and two smaller systems, to bracket a broad mass range across the considered redshift values (Table 1). In both mass bins, we choose a cool-core (CC) and a non-cool-core cluster (NCC), defined based on their pseudo-entropy profiles as described in Leccardi et al. (2010). This small sample gives a view of part of the expected cluster population planned to be investigated by the X-IFU. The objects are part of a larger set of 29 Lagrangian regions extracted from a parent cosmological DM-only simulation and re-simulated at higher resolution including baryons (see Bonafede et al. 2011). The parent cosmological volume is $1 h^{-1}$ Gpc per side and adopts a Λ -CDM cosmological model with $\Omega_{\text{M}} = 0.24$, $\Omega_{\text{b}} = 0.04$, $H_0 = 72 \text{ km s}^{-1} \text{ Mpc}^{-1}$ (i.e. $H_0 = h \times H_{100}$, where $h = 0.72$ and $H_{100} = 100 \text{ km s}^{-1} \text{ Mpc}^{-1}$), $\sigma_8 = 0.8$ and $n_s = 0.96$, consistent with WMAP-7 constraints given in Komatsu et al. (2011).

The magnification simulations were performed with a version of the tree-PM smoothed particle hydrodynamics (SPH) code GADGET-3 (Springel 2005), including an improved hydrodynamical scheme (Beck et al. 2016) and a variety of physical processes describing the evolution of the baryonic component (see Rasia et al. 2015, for more details). Briefly, these comprise metallicity-dependent radiative cooling (Wiersma et al. 2009), star formation and stellar feedback (thermal supernova feedback and galactic winds, see Springel & Hernquist 2003), cold and hot gas accretion onto super-massive black holes powering AGN thermal feedback (Steinborn et al. 2015; modelling the action of cold accretion Gaspari & Sdowski 2017), and metal enrichment (Tornatore et al. 2004, 2007)

¹ R_{500} is the radius including a density contrast of 500 times the critical density of the Universe, $\rho_c = 3H(z)^2/8\pi G$, at the given redshift z .

Table 1. Properties of the simulated clusters at different redshift values in their evolution.

Name		C1	C2	C3	C4
Type		CC	NCC	CC	NCC
$z = 0.105$	R_{500} (kpc h^{-1})	723	799	1027	1009
	M_{500} ($M_{\odot} h^{-1}$)	2.39×10^{14}	3.22×10^{14}	6.86×10^{14}	6.51×10^{14}
	T_{500} (keV)	3.22	4.20	6.47	6.36
$z = 0.5$	R_{500} (kpc h^{-1})	552	676	694	715
	M_{500} ($M_{\odot} h^{-1}$)	1.55×10^{14}	2.84×10^{14}	3.08×10^{14}	3.38×10^{14}
	T_{500} (keV)	2.88	4.06	4.58	4.84
$z = 1.0$	R_{500} (kpc h^{-1})	389	396	446	570
	M_{500} ($M_{\odot} h^{-1}$)	0.92×10^{14}	0.97×10^{14}	1.38×10^{14}	2.89×10^{14}
	T_{500} (keV)	2.41	2.47	3.12	4.41
$z = 1.48$	R_{500} (kpc h^{-1})	269	289	345	351
	M_{500} ($M_{\odot} h^{-1}$)	0.50×10^{14}	0.62×10^{14}	1.07×10^{14}	1.12×10^{14}
	T_{500} (keV)	1.76	2.10	3.07	3.08
$z = 2.0$	R_{500} (kpc h^{-1})	174	181	220	215
	M_{500} ($M_{\odot} h^{-1}$)	0.23×10^{14}	0.25×10^{14}	0.45×10^{14}	0.42×10^{14}
	T_{500} (keV)	1.19	1.55	2.24	1.86

from SN_{Ia}, SN_{cc}, and AGB stars. Specifically, we assumed the IMF by Chabrier (2003), the mass-dependent lifetimes by Padovani & Matteucci (1993) and stellar yields by Thielemann et al. (2003) for SN_{Ia}, Woosley & Weaver (1995) and Romano et al. (2010) for SN_{cc}, and Karakas (2010) for AGB stars.

In our model of chemical enrichment, we follow the production and evolution of 15 chemical species: H, He, C, Ca, O, N, Ne, Mg, S, Si, Fe, Na, Al, Ar, and Ni. These elements are the individual species traced in the simulations. Although these do not cover the full spectrum of interest (lacking e.g. Mn or Cr, which are important tracers of the enrichment as recently shown in Hitomi Collaboration 2017; Simionescu et al. 2018), the variety of abundances provides a good starting point for a meaningful study on the ICM and the demonstration of the X-IFU capabilities in this view. For every gas particle in the simulation, we traced the chemical composition and the fraction of each metal that is produced by the three enrichment sources (i.e. SN_{Ia}, SN_{cc}, AGB; see Biffi et al. 2017, 2018, for further detail). Each object is analysed at different redshifts, $z = 0.105, 0.5, 1, 1.48,$ and 2 , to assess the enrichment through time. Table 1 provides the characteristic radius, R_{500} , along with the mass, M_{500} , and the mass-weighted temperature, T_{500} , of the associated sphere of radius R_{500} for the entire cluster sample.

For each SPH particle, the output quantities provided by GADGET-3 are used as input for our simulation. These include the position of the particle, \underline{x} , its 3D velocity in the observer's frame, \underline{v} , its mass density, ρ , its mass, m , its temperature, T , and the individual masses of the 15 individual chemical species X, μ_X , tracked in the simulations. The gas density n of each SPH particle is obtained by dividing ρ by m . The mass of each element of atomic mass number A_X is converted into abundances Z_X , expressed in solar metallicity units assuming the solar fractions $Z_{\odot,X}$ from Anders & Grevesse (1989). Abundances can be therefore written as

$$Z_X = \frac{1}{Z_{\odot,X}} \times \frac{\mu_X}{\mu_H \times A_X}, \quad (1)$$

with μ_H the hydrogen mass of the particle.

3. End-to-end simulations

In this section, we describe in detail the set-up of the pipeline used for the synthetic X-IFU observations, as well as the physical assumptions made in the simulations.

3.1. Synthetic X-IFU observations

Simulations of the cluster data set are carried out using the X-IFU end-to-end (E2E) simulator SIXTE² (Wilms et al. 2014), which creates realistic X-IFU observations. SIXTE uses as an input a specific SIMPUT file (Schmid et al. 2013) containing either all the emission spectra of the particles or directly a photon list, with the time, coordinates on the sky and energy of the emitted photons. This second approach is preferred for our simulations, as it exerts a lower computational demand, induced in the former by the unparallelised random generation of photons currently implemented in SIXTE (an example of the first approach is given in Roncarelli et al. 2018). SIXTE outputs are generated not only considering the instrumental spatial and spectral responses, but also incorporate other features from the detectors such as their geometry, vignetting and internal particle background.

3.1.1. Photon list generation

Each simulated cluster comes as a list of SPH particles, which may emit X-ray photons. To generate the photon list used in the E2E simulation, the particle emission is modelled by a collisional diffuse thermal plasma using the APEC code (Smith et al. 2001). More specifically, the vvapec model on XSPEC (Arnaud 1996) is adopted, as it can be parametrised according to the particles physical properties listed above, notably the individual abundances of each element. The corresponding atomic database used for the emission model is derived from ATOMDB v3.0.9. For the galactic absorption, the wabs model (Morrison & McCammon 1983) is preferred for computational speed, although more accurate absorption models do exist (e.g. TBabs, Wilms et al. 2000). For all four clusters

² <http://www.sternwarte.uni-erlangen.de/research/sixte/>

we fixed the column density to $n_{\text{H}} = 0.03 \times 10^{22} \text{ cm}^{-3}$, which is a representative value for the latitudes at which most clusters shall be observed with X-IFU (Kalberla et al. 2005). Abundances are set to solar as per Anders & Grevesse (1989) and atomic cross-sections are taken as per Verner et al. (1996). The overall flux, F (in $\text{counts s}^{-1} \text{ cm}^{-2}$), of each particle is computed using the `vvapec` normalisation \mathcal{N} (emission-measure-weighted by the distance in units of cm^{-5}):

$$\mathcal{N} = \frac{10^{-14}}{4\pi[D_{\text{A}}(1+z)]^2} \int n_{\text{e}}n_{\text{p}}dV, \quad (2)$$

where D_{A} is the angular distance of the particle computed from its redshift z (derived from the speed of the particle and the cluster mean redshift), and V the particle volume. We considered a full ionisation of the intra-cluster gas with $n_{\text{e}} = 1.2n_{\text{p}}$ (n_{e} and n_{p} being the densities of electrons and protons, respectively). These emission spectra are considered as probability density distribution function and normalised accordingly over the instrumental energy bandpass (i.e. 0.2–12 keV). For a fixed exposure time Δt , photons are drawn from the afferent probability distribution following a Poisson statistic of parameter $F\Delta tA$, where A is the total mirror area (taken at 1.4 m^2 at 1 keV, energy dependence of the effective area is included later on in SIXTE via the ancillary response function – ARF as explained below). Each newly created photon is added to the photon list with the sky coordinates of its parent particle (right ascension and declination).

The output product of this stage is a “complete” photon list (with their true energy and position) at the entrance of the telescope. This list is computed once for each cluster, and contains a large number of simulated photons ($\geq 1 \text{ Ms}$). It is then sampled randomly by SIXTE to achieve smaller lists for more typical exposure times (e.g. 100 ks).

3.1.2. Observational setup

For each simulation, we consider an exposure time $\Delta t = 100 \text{ ks}$ over the entire X-IFU field-of-view. The complete photon list is used as input for the `xifupipeline` function of SIXTE, which samples the photon list accounting for the energy-dependence of the effective area to create the event list seen by the X-IFU detector over Δt . The pipeline accounts for the most up-to-date responses of the current baseline of the telescope (i.e. 15-row mirror modules corresponding to a mirror effective area of 1.4 m^2 at 1 keV^3) and for a hexagonal detector array of 3832 micro-calorimeter pixels, more specifically Large Pixel Array 2 (LPA2) pixel configuration, developed for the X-IFU and described in Smith et al. (2016). Pile-up, telescope point spread function, vignetting and detector geometry effects are also included as function of the pixels corresponding off-axis angles. Finally, we verified that given the low count rates of our clusters ($\leq 1 \text{ cts s}^{-1} \text{ pix}^{-1}$), pile-up and cross-talk over the observation can be neglected (see den Hartog et al. 2018; Peille et al. 2018).

For each cluster, we simulated enough pointings to map the cluster spatially up to at least R_{500} (as required in the current science objectives for the X-IFU). This translates, for local clusters, into at least seven pointings. The corresponding event lists are then merged during post-processing to obtain a single event file.

³ RMF: `athena_xifu_rmf_v20171107.rmf` | ARF: `athena_xifu_15row_onaxis_pitch249um_v20171107.arf`

Table 2. Parametrisation of the galactic foreground model used in the simulation with a `apec + phabs*apec` model.

Model	Parameter	Unit	Value
apec	T	keV	0.099
apec	Z		1
apec	z		0
apec	Norm	$\text{cts s}^{-1} \text{ amin}^{-2}$	1.7×10^{-6}
phabs	n_{H}	10^{22} cm^{-3}	0.018
apec	T	keV	0.225
apec	Z		1
apec	z		0
apec	Norm	$\text{cts s}^{-1} \text{ amin}^{-2}$	7.3×10^{-7}

3.2. Foreground and background components

In addition, we accounted for the contribution of different foreground and background sources to ensure more representative observations.

3.2.1. Astrophysical foreground

The foreground emission is caused by the X-ray emission of the local bubble in which the solar system is embedded and by the Milky Way hot gaseous content. This component can be modelled by the sum of a non-absorbed and absorbed thermal plasma emission as specified in McCammon et al. (2002) and parametrised as per (Lotti et al. 2014, see Table 2). An additional normalisation constant over the entire model is used for versatility purposes, resulting in a total foreground model reading as `constant*(apec + phabs*apec)` in XSPEC. This component is folded into SIXTE using a SIMPUT file.

3.2.2. Cosmic X-ray Background

The cosmic X-ray background (CXB) component is due to the contributions of AGNs, star forming galaxies and active stars along the line-of-sight (Lehmer et al. 2012). A fraction of these sources will be resolved by the instrument as a function of its spatial resolution, and will be excised from the observations. Given the requirement on the spatial resolution for *Athena*/X-IFU ($5''$), 80% of the total flux of these point sources in terms of the integral of their $\log(N)/\log(S)$ distribution should be resolved by the instrument (Moretti et al. 2003). For 100 ks exposure times, this translates into limiting fluxes of $\sim 3 \times 10^{-16} \text{ ergs s}^{-1} \text{ cm}^{-2}$ for the X-IFU. As the number of star forming sources is at least an order of magnitude lower at this flux, we only considered the AGN contribution in this study. The unresolved fraction of these point sources results in a diffuse background component, which we classically fitted using an absorbed power-law model during post-processing (McCammon et al. 2002).

We did not include AGN point sources in the inputs derived from the hydrodynamical simulations. Instead, to generate realistic CXB data, we drew a list of AGN sources with associated X-ray spectra by sampling the luminosity function of Hasinger et al. (2005) in the luminosity-redshift space, given the boundary conditions $L_{\text{X}} \geq 10^{42} \text{ erg s}^{-1}$ unabsorbed 0.5–2 keV rest-frame luminosity, $0 < z < 5$ and the size of the cosmological volume encompassed within a field-of-view. Each source is associated with a spectral energy distribution following templates described in Gilli et al. (2007), according to a distribution

can no longer be detected (lines outside the energy bandpass) and rarer elements (e.g. Ne, Na, and Al) have large uncertainties due to the low S/N of the observations. Ni also tends to be underestimated (mostly in the outskirts) likely due to the low S/N of the line with respect to the high-energy background. This calls for better-adapted exposure strategies to optimise the results for distant objects and further investigate the chemical enrichment across cosmic time.

7. Summary and discussion

In this paper, we have addressed the feasibility of constraining the chemical enrichment of the Universe, which will be one of the key science objectives and a main driver of the performances of the future mission *Athena*. Notably, we investigated and quantified the capabilities of the X-IFU in accurately recovering metal abundances of the ICM across cosmic time. To this end, we developed a full end-to-end pipeline, which creates synthetic X-IFU observations using the instrument simulator SIXTE. We used as input of this pipeline a sample of four clusters generated using state-of-the-art cosmological simulations presented in [Rasia et al. \(2015\)](#) and [Biffi et al. \(2017\)](#) to create realistic event lists. All the relevant instrumental effects such as the convolution of spectra with the instrument spatial and spectral responses, realistic sources of foreground and background, and detector geometry were also included to obtain observations as realistic as possible.

The sample of four clusters was simulated at five different redshift values, for a fixed exposure time of 100 ks in order to achieve abundance measurements out to R_{500} . The accuracy of the pipeline was quantified by comparing our synthetic observations to weighted inputs quantities (e.g. spectroscopic temperature, emission-measure-weighted abundances). We find that a straightforward approach of a broad-band fit created systematic biases above 10% in a number of physical parameters. Rather a multi-band energy fitting procedure ensured more accurate recovery by optimising the extraction of the several chemical abundances and other physical parameters of interest (notably temperature). After post-processing, distributions are accurately recovered (almost always within the 3σ of the measurement error) with little to no systematic biases (of the order of a 5%, see Sect. 4.6) found mainly between the low-mass element abundances (e.g. O, Si) and the normalisation. The comparison of the relative distribution between outputs and inputs with respect to the XSPEC statistical error also showed reduced chi-squared values χ^2_{red} close to 1 when a small fraction of outlier regions ($\leq 5\%$) is removed indicating a good accuracy in the fits. Remaining errors and biases can be linked to correlation between parameters (notably the normalisation), to the choice of the input weighting scheme, and to mixing effects along the line-of-sight in view of the single plasma temperature model used here. We also find that most of these errors decrease when statistics are strongly increased (biases below 2% at 1 Ms for the same spatial regions), suggesting that these effects may simply be related to statistics (see Appendix B). Studies conducted by decreasing the statistics of the runs (typically by decreasing the S/N of the regions to 50 or 100) provided equally encouraging results. Despite larger statistical errors (up to 10% higher), the main parameters (temperature, redshift, O, Si, or Fe abundance) were accurately recovered. Some fainter lines (e.g. Na, N, or Al) become however very difficult to constrain in this case.

For local clusters ($z \sim 0.1$), we demonstrate the power of the X-IFU in accurately recovering spatially resolved parameter maps, along with abundance profiles (Sect. 5.2) and abundance ratios (Sect. 5.3). The study was then extended at different red-

shift values, up to $z \sim 2$. By probing the chemical enrichment for very distant clusters and despite the lack of an optimised observation strategy (i.e. non-optimised exposure time), we also show the power of the X-IFU in investigating the ICM properties and the chemical enrichment of the distant Universe.

The binning and fitting procedures used here comprise “classical” approaches to X-ray data analysis, using S/N binning and fits through instrumental response matrices in XSPEC. Despite our efforts, the fitting procedures remain slightly biased ($\leq 5\%$) and small changes in the fitting approach can impact the overall results of the simulation (of the order of a few %). More accurate results may be achieved using, for example, Monte-Carlo (MC) fitting approaches, but unfortunately remain computationally cumbersome to be used on our large set of spatial regions. More optimised binning techniques could also be investigated for future applications ([Kaastra & Bleeker 2016](#)). The access to high-resolution spectra will provide new proxies to estimate quantities such as the temperature by using for example, line-ratio techniques. Eventually, hyper-spectral methods (e.g. blind source separation algorithms) or machine-learning-based fitting techniques (see, e.g. [Ichinohe et al. 2018](#)) could open new perspectives for the post-processing of high-resolution X-ray spectra. We would like to emphasise that, even though not applicable in our simulation case, the expected level of spectral resolution of the X-IFU will challenge our current knowledge accuracy of the spectral lines (centroid energies and intrinsic widths). This is critical to allow a meaningful interpretation of the results (as demonstrated in [Hitomi Collaboration 2018d](#), for line ratios) and to disentangle fine spectroscopic effects (such as resonant scattering, [Hitomi Collaboration 2018b](#)). This emphasises the need for dedicated tools able to process and analyse future X-IFU high-resolution spectroscopy data-cube. In this regard, the *Athena* mission will certainly benefit from the advances expected in processing tools, fitting methods and atomic databases, from the future XRISM mission ([Ishisaki et al. 2018](#)).

Not only do these E2E simulations allow us to explore the capabilities of the future X-IFU instrument, but they also give crucial information on the effect of instrumental parameters in science observations. In this study for instance, the spectral shape of all the foreground and background components were assumed to be perfectly known. For the more local and massive clusters however, the field-of-view of the X-IFU will easily be encompassed within the angular extension of R_{500} . Cluster emission-free regions might thus be unavailable for local background calibration. The spectral resolution of the X-IFU will help mitigate this effect, by allowing us to disentangle various components through the characteristics of their spectral energy distribution. The instrument background may also contaminate the observation of faint sources, as the level of precision to which X-IFU is expected to perform requires its accurate and reproducible knowledge in flight. This may be achieved through, for example, in-flight cross-correlation with the WFI or the X-IFU cryogenic anti-coincidence detector ([Cucchetti et al. 2018](#)). Future developments could take advantage of this simulation pipeline to test other realistic instrumental effects (e.g. stray-light for galaxy cluster outskirts observations). More detailed studies of the abundance ratios recovered here will also be at the centre of a forthcoming study to highlight the capabilities of the X-IFU in constraining the ICM chemical enrichment, and notably to disentangle between the contributions of the various mechanisms of chemical enrichment (e.g. SN, AGB) throughout cosmic time.

Our study underlines the revolutionary capabilities brought by the X-IFU in future X-ray spectroscopy. With typical routine observations, the X-IFU will drastically change our

understanding of ICM mechanisms and provide a quantum leap forward in X-ray astronomy.

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Appendix A: Spatial binning algorithm comparison

Spatial binning of the data is used to increase the S/N of the considered regions, to have higher significance spectra. This is of particular interest whenever fine structures (e.g. line ratios, line doublets, absorption features) need to be observed in a spectrum. Multiple methods can be used to bin spatially, in this study two of them were considered:

- `contbin` tool developed by J. Sanders (Sanders 2006). The `contbin` scheme was run for a constrain fill value of two, which represents the maximum ratio length/width of a region.
- Voronoï tessellation, as defined by (Cappellari & Copin 2003)

Both these methods were tested on the same surface brightness map to estimate their performances. Various criteria were compared, such as their ability to reproduce spatial features of the cluster or the mean S/N ratio of the regions created. Figure A.1 shows a comparison of the cluster 2 regions at $z = 0.105$ without any foreground/background component (to investigate purely the binning effects) computed using either of these methods. Visually, we notice that `contbin` provides very similar results to Voronoï whenever the aspect ratio of the region is constrained to $C_{\text{fill}} = 1$ (Fig. A.1, right). When a slightly higher value of the aspect ratio is allowed (Fig. A.1, bottom left), we notice that `contbin` is able to reproduce more accurately the radial contours of the cluster, notably the cold arc visible in the south-east corner of input data or the hot bubble rising west of the cluster (Fig. A.1, top left). Further, no difference is found on the average S/N of the regions, which is always above the required level. Finally, the methods give very close results in terms of number of regions (85 for Voronoï vs 87 for `contbin` for a S/N of 300), thus being equivalent computationally. For our purposes, `contbin` tool provides a more suitable binning algorithm than Voronoï. More ample tests also show no significant difference in the recovery of the physical parameters between both techniques.

Appendix B: Validation of the simulation pipeline

Test hypothesis

The accuracy of the simulation pipeline needs to be verified using the cluster inputs provided. These inputs are 3D cubes of data, which needs to be projected along the line-of-sight of the instrument to be compared to the outputs of the end-to-end simulations. Ideally, this projection should be deterministic and give an unequivocal results. However, since the parameter distribution along the line-of-sight cannot be perfectly integrated (we only measure discrete number of counts, affected by statistics and background sources), multiple schemes exist to compare inputs and outputs depending on the physical quantity we wish to compare. Among those, the most widely used include:

- Emission-weighted projection, using the product $\rho^2 \sqrt{T}$ of each element along the line-of-sight.
- Mass-weighted projection, using the mass of each element
- Emission-measure-weighted projection, using the emission-measure of a given line derived from Eq. (2)
- Spectroscopic schemes, as defined in (Mazzotta et al. 2004)

The accuracy of the simulated measurements was first tested by taking as estimator the relative error distribution of the output map, using emission-weighted-input maps as proxy for the input parameters. Being each region much larger than the telescope PSF they are considered independent on a strict statistical term. The relative error is assumed to follow a Gaussian

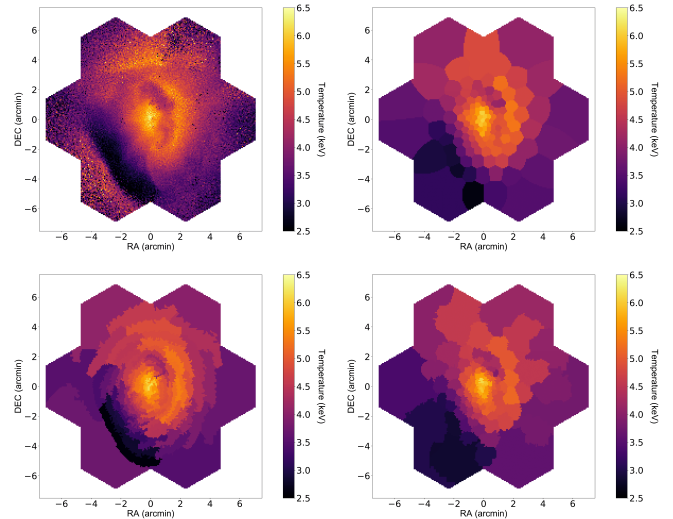


Fig. A.1. Spatial binning scheme comparison for cluster 4 spectroscopic temperature map (keV) without background, at redshift $z = 0.105$. *Top left:* unbinned raw input map from the hydrodynamical simulation. *Top right:* voronoï tessellation map using the algorithms described in (Cappellari & Copin 2003). *Bottom left:* contour map using `contbin` tool (Sanders 2006) with an aspect ratio constraint, $C_{\text{fill}} = 2$. *Bottom right:* same as *bottom left*, with $C_{\text{fill}} = 1$.

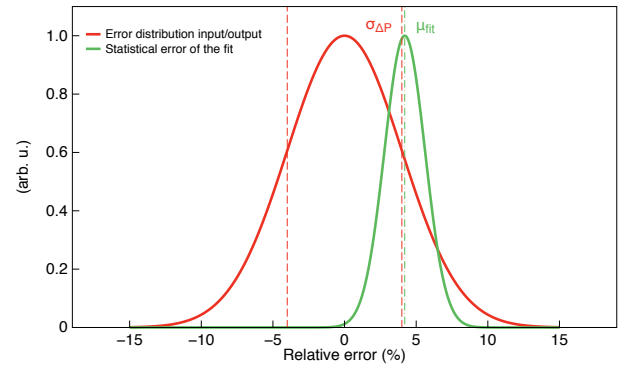


Fig. B.1. Comparison between the relative error distribution of the parameters (red) and the fitting error distribution (green – symmetric but not shown here for clarity). Ideally, if no biases are present relative error distribution should be centred and its standard deviation $\sigma_{\Delta P}$ should be very close to the mean value of the fitting error distribution μ_{fit} .

distribution given the sufficiently high number of regions considered (≥ 80), with a mean $\mu_{\Delta P} = 0$ (if no biases are present) and a standard deviation of $\sigma_{\Delta P}$, which indicates the total error on the parameter. The fitting error returned by XSPEC should also be Gaussian, and centred on given a value μ_{fit} which depends on the exposure time and the emission model parameters. For an accurate measurement, the value of μ_{fit} should be comparable to $\sigma_{\Delta P}$ do all parameter (Fig. B.1). A second test can also be performed using as estimator the ratio between the relative difference and the XSPEC error σ_{fit} for each region, i.e. $\chi_j = (P_{\text{fit},j} - P_{\text{in},j}) / \sigma_{\text{fit},j}$ and the corresponding reduced chi-square. Using the emission-measure-weighted input and the output distributions, let us take as null hypothesis (H_0): “The measurements obtained with the pipeline are consistent with the statistical errors for a given exposure time” and (H_1): “The measurements are biased” with a threshold $p_\alpha = 5\%$ (i.e. 97.5% of the Gaussian distribution, or $\sim 2.5\sigma$).

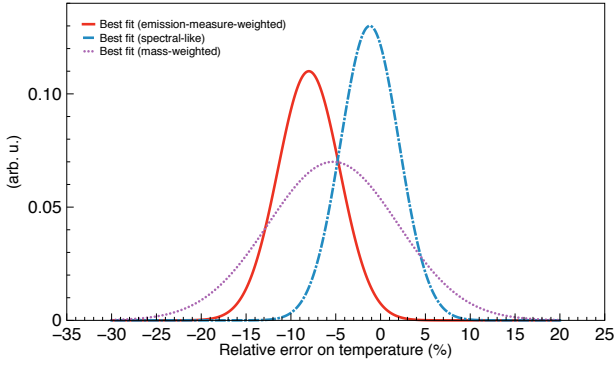


Fig. B.2. Gaussian best fits of the normalised relative error distribution on the measured temperature for different input map weighting scheme for cluster C4. The red solid curve indicates the emission-weighted best fit ($\mu_{\Delta T} = -8.0\%$, $\sigma_{\Delta T} = 3.4\%$). The blue dash-dotted line indicates the spectroscopic temperature best fit ($\mu_{\Delta T} = -1.2\%$, $\sigma_{\Delta T} = 2.2\%$), while the dotted violet line indicates the best fit for a mass-weighted input ($\mu_{\Delta T} = -5.2\%$, $\sigma_{\Delta T} = 7.2\%$).

H_0 : (H_0) can be rejected if the value of $\sigma_{\Delta P}$ is outside $\mu_{\text{fit}} \pm 2.5\sigma_{\text{fit}}$ for all the parameters. This can also be seen on the χ_j maps, if the number of regions with $|\chi_j| \geq 2.5$ is high. For all runs, the value of the error dispersion is always within this threshold and the number of outliers regions is small (see also Appendix C and Table 3), indicating that the (H_0) is valid and errors are consistent with the corresponding statistics.

H_1 : (H_1) can be rejected if $\mu_{\Delta P} \sim 0$, within ~ 2.5 times the mean standard error of the distribution (due to the finite size of the sample). The samples are generally composed of $N_{\text{reg}} \sim 80$ regions and the standard error on $\mu_{\Delta P}$ is given by $\sigma_{\Delta P} / \sqrt{N_{\text{reg}}}$. Under these assumptions, clear biases are visible in the reconstructed emission-weighted and mass-weighted temperature maps, which presented a systematic underestimation of ~ 5 – 10% (Fig. B.2). This bias is explained by mixing effects along the line-of-sight and complexity to disentangle multi-temperature plasma with a single plasma model (Mazzotta et al. 2004). It can be reduced using spectroscopic temperature maps (Fig. B.2). The use of the broad-band fit induced in fact many other visible biases, notably between abundances (O, Si, and Fe) and temperature (Fig. B.3, upper panel). The use of multi-band fitting (detailed in Sect. 4.6) significantly reduces these biases (Fig. B.3, lower panel) within the statistical variations of the parameters. Despite this improvement, small correlations are visible between abundances and temperature ($\sim 1\%$) and between the normalisation and all other parameters ($\sim 5\%$). Efforts to reduce this bias were conducted by fixing or releasing various fitting parameters, without significant success. A clear rejection of the null hypothesis (H_1) cannot be performed, although results suggest that small residual biases and correlations remain in the current fitting procedure, mainly on normalisation.

Influence of the projection scheme

Depending on the weighting schemes (mass-weighted or emission-measure-weighted) different error distributions of the same quantities can be obtained. These discrepancies add a further complexity to evaluate any potential bias in the pipeline (Fig. B.2).

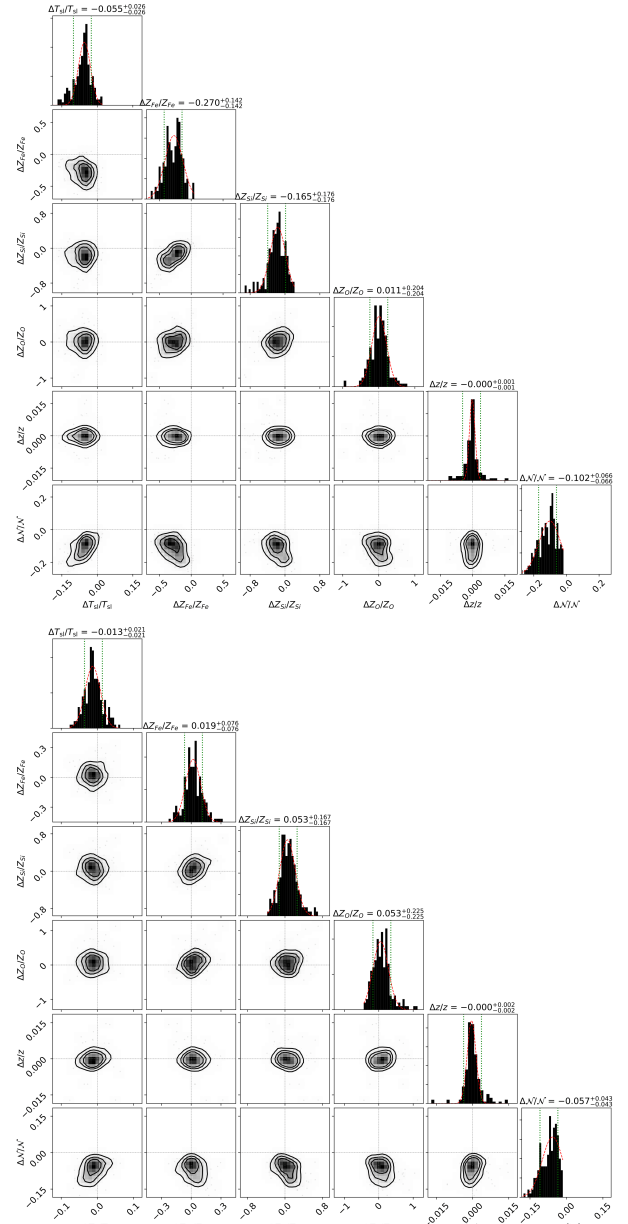


Fig. B.3. Corner plots of the relative error on parameter $(X_{\text{fit}} - X_{\text{inp}})/X_{\text{inp}} = \Delta X/X_{\text{inp}}$ as function of parameter $(Y_{\text{fit}} - Y_{\text{inp}})/Y_{\text{inp}} = \Delta Y/Y_{\text{inp}}$, for the spectroscopic temperature, T_{sl} , oxygen, silicon, and iron abundance, redshift, z , and the normalisation N , for cluster C4. The diagonal panels are the corresponding relative error distribution, where the red solid line indicates the Gaussian best fit of the distribution (parameters $\mu_{\Delta P}$, $\sigma_{\Delta P}$ given above) and the dotted line is the value of $\mu_{\Delta P} \pm \mu_{\text{fit}}$. *Top*: Broad-band fit. *Bottom*: multi-band fit (Sect. 4.6) considering a spectroscopic temperature.

Accuracy in terms of probability distributions

The previous test is only valid whenever the error distributions are assumed to be Gaussian, which is unfortunately not always the case (slight deviations from Gaussian behaviour are observed). If so, the accuracy of our method needs to be tested in the sense of the statistical distributions performing for instance a Kolmogorov–Smirnov (KS) test over the output and input distributions. The KS test compares the probability p_{KS} for two random variables to be drawn from the same data set (i.e., same

probability density). Using both the input and output distribution, this test showed high p_{KS} values (between 0.6 and 1 for the different parameters), which gives strong hints that the output distribution indeed matches the input. Statistically speaking, no real conclusion can be achieved with a single realization of the observation. To fully validate the pipeline, a large number of observations of the same cluster (either along the same line-of-sight or by taking multiple lines-of-sight) would be needed to perform a meaningful comparison using a KS method.

Unfortunately, the duration of one full simulation of a cluster is of the order of a day, making it computationally cumbersome to carry out this test. For simplicity, a very high exposure time simulation of these extended sources were carried out instead. Although beyond the scope of this paper, such observations (≥ 1 Ms) with the same binning regions decrease most of the biases below 2% and create distributions which much more alike ($p_{KS} \approx 0.8/1$), suggesting that the residual errors are in part related to statistics and to the fitting scheme.