

# Spatially resolved spectroscopy of Cassiopeia A with MECS on board *BeppoSAX*

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**Abstract.** We have performed the first detailed spatially resolved spectroscopy of Cas A in the 1.6–10 keV energy range, using data taken with the MECS spectrometer on board the *BeppoSAX* Observatory. The well calibrated point spread function in the central region of the MECS allowed us to perform a spatial deconvolution of the data at full energy resolution. We eventually generated a set of spectra, covering a region of  $\sim 3'$  radius around the centre of Cas A. The results obtained by fitting these spectra using a non-equilibrium ionisation plasma model and a power law, improve our knowledge about chemical and physical parameters of the Cas A SuperNova Remnant: (i) a single thermal component is sufficient to fit all the spectra; (ii)  $kT$  is rather uniformly distributed with a minimum in the east and a maximum in the west, and no evidence is found for high  $kT$  expected from the interaction of the main shock with the ISM; (iii) from the distribution of the values of the ionisation parameter  $n_e t$  we infer the presence of two distinct components: the first ( $a$ ) with  $n_e$  in the range  $1\text{--}10\text{ cm}^{-3}$ , the second ( $b$ ) with values ten times higher; if we associate component  $a$  to the CSM and component  $b$  to the ejecta, the mass ratio  $M(a)/M(b) \leq 1/10$  indicates a progenitor star that lost only a small fraction of the envelope during its pre-SN life. In this hypothesis the distribution of component  $b$  across the remnant suggests that the explosion was not spherically symmetric; (iv) the distribution of abundances indicates that we are detecting a CSM component with almost solar composition, and an ejecta component enriched in heavier elements. Abundances found for  $\alpha$ -elements are consistent with the current view that Cas A was produced by the explosion of a massive star. A low Fe overabundance can be an indication that at the moment of the explosion the mass-cut was rather high, locking most of the produced Ni<sup>56</sup> into the stellar remnant.

**Key words.** ISM: SuperNova remnants – X-rays: ISM – methods: data analysis – techniques: image processing, spectroscopic – IMS: individual: Cas A

## 1. Introduction

Cassiopeia A (Cas A), the youngest known SuperNova Remnant (SNR) in our galaxy, is an object investigated at several wavelengths. It has the appearance of a broken shell with average radius of  $\sim 1.7'$  (Holt et al. 1994; Vink et al. 1996) around its expansion center that Reed et al. (1995) found located at  $RA_{2000} = 23^{\text{h}}23^{\text{m}}26^{\text{s}}.6$ ,  $DEC_{2000} = 58^{\circ}49'01''$ . Detailed observations at optical wavelengths show a complex morphology, not spherically symmetric, with two northwest (NW) and southeast (SE) emission regions moving outward from the remnant center in accordance with their relative average Doppler motion (Anderson et al. 1991; Reed et al. 1995; Lawrence et al. 1995; Fensen & Gunderson et al. 1996). A similar morphological complexity comes out in the X-ray band

as observed by ASCA at the moderate spatial resolution of  $3'$  (Holt et al. 1994; Fujimoto et al. 1995; Vink et al. 1996; Hwang & Gotthelf 1997). *BeppoSAX* (Boella et al. 1997a) observed the source giving interesting results either in the spectral and in the spatial analysis. In particular, the maps obtained in the energy range 1.6–10 keV (Maccarone et al. 1998; Vink et al. 1999) with the Medium Energy Concentrator Spectrometer MECS instrument (Boella et al. 1997b) show that the spatial distributions of Si, S, Ar, and Ca emission lines do not differ very much each other. In addition, the presence of Fe seems to extend further east than for the other elements. Such results have been confirmed recently by the Chandra X-ray observatory that, with its imaging capability of  $\simeq 0.5''$ , provided well detailed maps of the remnant and discovered a point-like source near its center (Hughes et al. 2000; Chakrabarty et al. 2000; Hwang et al. 2000).

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The broad band *BeppoSAX* spectrum of the whole source (Favata et al. 1997) presents a high energy component well modeled by a power law of spectral index 2.95. This component has been localized by Vink et al. (1999) mainly in the southern and western regions. Indeed the MECS spectra of these regions are somewhat harder than the average and almost featureless above 4 keV.

In this paper we present results about the spatially resolved spectroscopy of Cas A in the MECS energy range 1.6–10 keV, at the resolution of  $1.5'$ . Data set and data reduction are presented in Sect. 2 together with a brief description of the software environment. Image and spectral analyses are described in Sects. 3 and 4, respectively. Results are discussed in Sect. 5 while conclusions are presented in Sect. 6. The treatment of errors is in Appendix A.

## 2. Observations and data reduction

The Medium Energy Concentrator Spectrometer MECS is one of the narrow field instruments on board *BeppoSAX*. The MECS consists of three units (ME), each composed of a grazing incidence mirror unit with a position sensitive gas scintillation proportional counter located at the focal plane. The MECS field of view covers a radius of  $28'$  with position resolution of  $\sim 1.5'$ ; inside the central circular window of  $\sim 10'$  (delimited by a strongback structure) all the detector parameters are well defined.

The three ME units were all together active till May 6, 1997, when ME1 unit was switched off, due to a failure in the high voltage supply of the gas cell unit. ME2 and ME3 units continued to work normally.

The on-axis Point Spread Function (PSF) of the MECS can be modeled as sum of two components: a Gaussian, and a generalized Lorentzian. For each ME unit, at a given energy and at a given distance  $r$  from the detector center, the complete analytical expression for the on-axis PSF is (Boella et al. 1997b):

$$PSF(r, E, ME) = \frac{1}{2\pi \left[ R\sigma^2 + \frac{r_l^2}{2(m-1)} \right]} \times \left\{ R \exp\left(-\frac{r^2}{2\sigma^2}\right) + \left[ 1 + \left(\frac{r}{r_l}\right)^2 \right]^{-m} \right\}$$

where  $R$ ,  $\sigma$ ,  $r_l$  and  $m$  depend on the ME unit and are algebraic functions of the energy. The dependence from the unit can be averaged in the formulation: this allows us to consider a single PSF at each given energy. Moreover, the on-axis PSF can be considered valid within a radius of  $12'$  from the detector center (Molendi et al. 1997; Chiappetti et al. 1998).

Table 1 lists the data set reference codes used for the analysis presented in this paper. MECS data were reduced by using a standard procedure<sup>1</sup> with appropriate selection criteria so to avoid the South Atlantic Anomaly and to

**Table 1.** The Cas A data set used in this paper

Obs. Prog.	Observ. Date	Observ. Period	Archival Code	Exposure (s)	ME units
SVP	06aug96	743	30011002	27574	1, 2, 3
SVP	07aug96	745	30011001	26639	1, 2, 3
SVP	11sep96	899	30011003	16629	1, 2, 3
SVP	11sep96	901	30011003	16595	1, 2, 3
AO1	26nov97	2990	30159001	40789	2, 3

obtain the maximum rejection of background particles. A further selection was applied to take into account only data related to the orbital periods in which the  $Z$  star tracker (aligned with the narrow field instruments) was in use; in this case the attitude reconstruction achieves an uncertainty of the order of  $30''$  inside the strongback region. Under these conditions, the total exposure time for the MECS was  $\sim 128000$  s, the net exposure time for each observation is listed in Table 1.

Data reduction has been carried out using the *SAXDAS* v.1.2.0 package, developed under the *FTOOLS* environment. Image analysis and spectra generation were carried out under *MIDAS-97NOV* environment<sup>2</sup>; the context *saxmecs*<sup>3</sup> has been used to convert data in *MIDAS* format, to merge MECS observations in images, to generate the PSF as function of energy, and to accumulate the final spectra.

The adopted deconvolution procedure is based on the Lucy's deconvolution method (Lucy 1974), as implemented in *MIDAS-97NOV*, and it can be fruitfully applied on sources presenting an extended and structured morphology. With respect to our previous analysis (Vink et al. 1999), we reconsidered the deconvolution parameters: in particular, we rebinned data to  $24''$  per pixel and performed deconvolution at each energy channel with the monochromatic PSF computed for that channel, as detailed in Sect. 3.

The spectral analysis has been performed with the *XSPEC* v.9.00 software package.

## 3. Image analysis

To enhance the morphology of the Cas A spatial emission in the energy range 1.6–10 keV at the resolution of the MECS Pulse Invariant (PI) channel, we adopted the following procedure: for each PI channel in the chosen energy range (PI channel from 32 to 214) we have (i) selected data from each observation, (ii) converted data in  $X, Y$  sky pixels images, (iii) merged and rebinned images to improve statistics, (iv) generated the appropriate PSF. We then deconvolved each monochromatic rebinned image with its

<sup>2</sup> <http://www.eso.org/midas>

<sup>3</sup> <http://www.ifcai.pa.cnr.it/~sax/saxmecs/saxmecs.html>

<sup>1</sup> <http://www.sdc.asi.it/software/cookbook>

proper monochromatic PSF for a suitable number of iterations.

Rather large inaccuracies are present in the aspect solution of our original Cas A data set, particularly in the last observation (AO1) made after a gyroscope failure. To overcome the problem, present whenever we need to merge spatial data coming from different observations, we used a procedure similar to that adopted in Vink et al. (1999) and detailed in Maccarone & Mineo (1999)<sup>4</sup>.

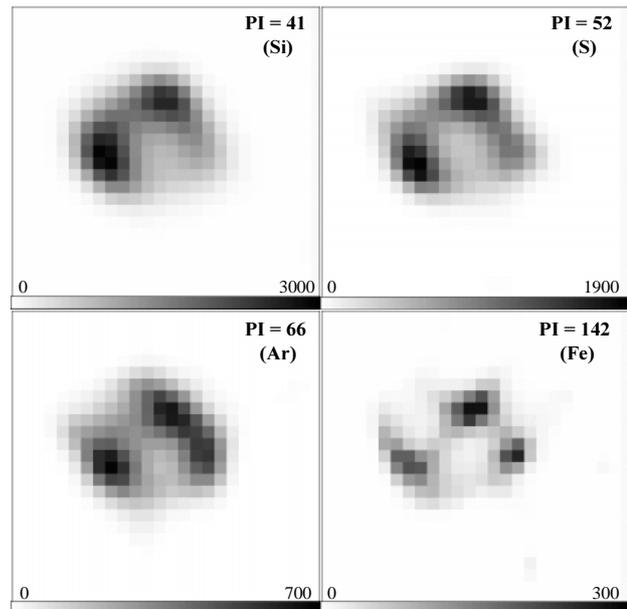
### 3.1. The deconvolution algorithm

To be successfully applied, the Lucy's deconvolution algorithm needs to be properly tailored taking into account the PSF characteristics, the spatial region of Cas A we are interested in, the statistical significance level we are dealing with, and the number of iterations necessary to reach convergence.

The spatial region of Cas A we are interested in is all inside the inner strongback window and it is centered very near to the detector center. Therefore, the radial description of the on-axis PSF can be successfully used in the deconvolution process. We applied the deconvolution algorithm to each PI channel within the chosen energy range by using the monochromatic PSF corresponding to that channel. The total number of counts within the interested region is about  $3.7 \cdot 10^6$  distributed over all the PI channels with a maximum of  $\sim 1.5 \cdot 10^5$  and a minimum of  $\sim 300$  counts.

To improve statistics, we compacted both PSF and observational data images by a factor of 3. This corresponds to rebin their stepsize from  $8''/\text{pixel}$  to  $24''/\text{pixel}$ , and to obtain a well-deconvolved spatial region of  $\sim 10'$  radius.

To define what number of iterations must be used in the deconvolution algorithm, we adopted a convergence criterion based on the relative change between images produced by two consecutive iterations. The iterations can be stopped when a given level of convergence for the maximum relative change is reached. From previous checks (Maccarone 1998; Vink et al. 1999), we noted that a maximum of 30 iterations is sufficient to enhance the morphology of the Cas A emission regions at higher energies (further iterations produce small corrections that tend to match the statistical fluctuations) while, at lower energies a greater number of iterations (some hundreds) is necessary to reach convergence. In the case of the Cas A data images analyzed here with stepsize of  $24''/\text{pixel}$  at the resolution of the MECS PI channel, a number of 200 iterations is sufficient to reach convergence at lower energies. To automatize the full deconvolution procedure we then choose to apply 200 iterations for any energy channel, without any risk of compromising the results for the higher energy channels. This choice was also supported by the comparison (evaluated via a standard  $\chi^2(i)$  function at each  $i$ th iteration) between the input image and the



**Fig. 1.** Deconvolution of the images related to PI channels 41, 52, 66, and 142, corresponding to  $\sim 1.89$ , 2.41, 3.06, and 6.66 keV, respectively

convolution of the  $i$ th deconvolved image with its proper PSF (Maccarone & Mineo 1999).

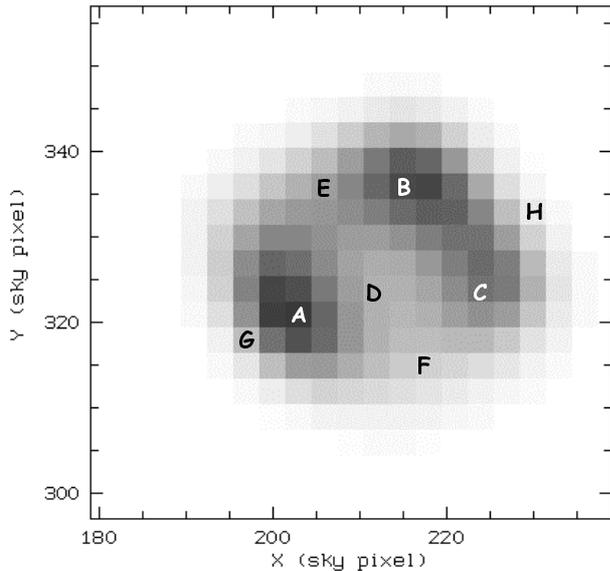
Figure 1 shows the results after 200 iterations on the images at the energy of Si, S, Ar, and Fe lines. The four monochromatic images present the same morphological structure already illustrated in Fig. 3, left panels of our previous paper (Vink et al. 1999), although the setting of the deconvolution parameters are slightly different:  $3 \times 3$  compacted pixels and full spectral resolution in this paper,  $2 \times 2$  compacted pixels and wide energy bands in our previous paper. From the image analysis point of view, the setting used here gives results in perfect agreement with those obtained in our previous paper; at the same time this setting allows us to perform a better spatially resolved spectroscopy at  $\sim 1.5'$  resolution level.

## 4. Spectral analysis

The set of Cas A images obtained from the deconvolution procedure has been organized in a position-energy cube  $(X, Y, E)$  from which we have accumulated spectra relative to each  $X, Y$  pixel in the spatial region of interest. To work with a good statistical signal we have only considered spectra with total counts greater than 800, reducing to 182 the final number of spatially resolved spectra to be analyzed. The window on which we have performed the spatially resolved spectroscopy then covers a region of  $\sim 3'$  radius around the Cas A expansion center. In sky pixels, the coordinates of this center are  $X = 209, Y = 327$ . Figure 2 shows the map of the total counts of the 182 selected pixels.

Statistical errors on the counts have been increased to take into account the systematics of the deconvolution

<sup>4</sup> <http://www.ifcai.pa.cnr.it/~cettina/papers/CasA-report2.ps.gz>



**Fig. 2.** X, Y projection of the Cas A data cube. The total counts range from 800 to  $\sim 62\,000$  counts/pixel. Labels refer to spectra shown in Fig. 3

procedure (see Appendix A); spectra have been rebinned to have at least 20 counts per energy bin.

Cas A is a strong source and the MECS background subtraction is not a relevant problem even if the deconvolution procedure acts on the spatial distribution of the background, moving counts from the external regions toward the center. We evaluated the background level in the worst case for which all counts within the well-deconvolved region ( $10'$ ) are placed under the source. Under the simplified hypothesis of an uniform background, the level is of the order of  $10^{-4}$  counts/s/pixel. We then ignored the background subtraction in our spectral analysis.

In the fitting procedure we used the MECS response matrix relative to an on-axis point source with an extraction circle of  $10'$  radius for all spectra, while the deconvolution was applied using a PSF defined in a squared area: this difference in geometry introduces an under-estimation of the effective area less than 1% in the spectral analysis.

Figure 3 shows a selection of spectra extracted at the positions marked with the corresponding letter in Fig. 2. It is evident that Cas A emission has not uniform spectral characteristics. In particular, as suggested by Vink et al. (1999) and Hughes et al. (2000), two main emission mechanisms can be identified: a thermal process, underlined by the presence of emission lines and coincident with the broken shell (Figs. 3-A, B, C, E, G), and an emission process characterized by the absence of strong lines (Figs. 3-D, F, H) and essentially located in the central region between the main peaks (Fig. 2).

To fit all spectra, we properly implemented in the fitting code a model of emission from a non-equilibrium ionization (NEI) plasma. The low energy absorption  $N_{\text{H}}$  has been constrained within the range  $0.9\text{--}1.7 \cdot 10^{22} \text{ cm}^{-2}$ , slightly wider than the range found by Keohane et al. (1996). The abundances of Si, S, Ar, Ca, Fe and Ni have

been left as free parameters and allowed to vary in the range 0.02–20 (with respect to solar). Moreover we added to the thermal model a power law to take into account the 20–80 keV emission found by Favata et al. (1997). The parameters of the power law are not free in the fitting procedure. In fact, none of the MECS deconvolved spectra has enough statistical significance to justify the presence of a second spectral component (e.g., the power law) superimposed to the thermal emission model.

The spectral index has been fixed at the value of 2.95 as quoted in Favata et al. (1997); the normalization has been computed by assuming that the whole remnant contributes to the high energy component and that the emissivity of this component is proportional to the MECS flux detected in each pixel. In the framework of the present analysis, we realized that the hard tail cannot be located in the two regions, south and west, suggested in our previous analysis (Vink et al. 1999). In fact, by fitting with a single power law in the 4–10 keV range the continuum emission of the spectra of these regions (covering  $\simeq 50\%$  of the remnant), and extrapolating the power laws in the high energy range, we found that the extrapolated flux is insufficient by a factor  $\sim 2$  to justify the high energy detection.

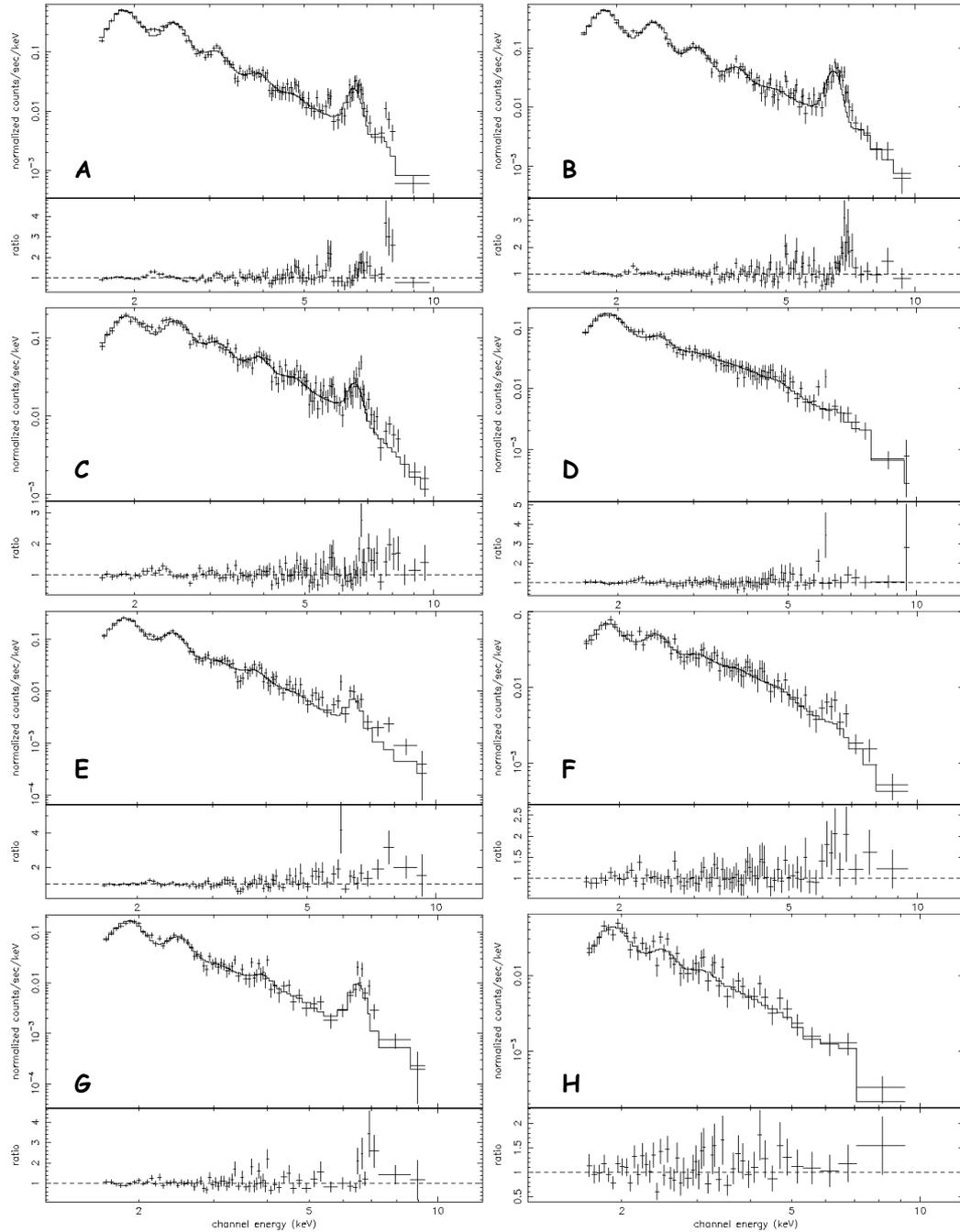
While the paper was in the refereeing process, results from XMM-Newton observations of Cas A became available (Bleeker et al. 2000), fully confirming our assumption. Indeed they state: “the hard X-ray image and the hardness ratio indicate that” the 8.1–15 keV “flux does not predominate in a few localized regions, but pervades the whole remnant in a distribution similar to the softer thermal component”.

Most of the spectra are well fitted by the composite model: the  $\chi^2$  values are within the expected range and no systematic residuals are observed in any of the spectra (see Fig. 3).

## 5. Results and discussion

Results obtained fitting the 182 spectra with a NEI plasma model are listed in Table 4. They are also shown in Figs. 4 and 5, where we present an histogram of the frequency distribution (left panels) and a map of the spatial distribution of the best fit values for each variable. In addition, we present in the right panel of Figs. 4 and 5 a frequency histogram of the same data used in the left panel, but where each single determination is transformed into a normalized Gaussian with central value and  $\sigma$  as derived from the fits.

We note that in previous studies analyzing data from a variety of different instruments including *BeppoSAX* (Hill et al. 1975, from a rocket-borne proportional counter; Davison et al. 1976, experiment C on *Ariel-5*; Pravdo et al. 1976, proportional counter on *OSO-8*; Becker et al. 1979, SSS on *Einstein*; Jansen et al. 1988, EXOSAT; Holt et al. 1994, ASCA; Vink et al. 1996, ASCA; Favata et al. 1997, *BeppoSAX*), reasonable fits of the spatially unresolved spectrum of Cas A were obtained using two



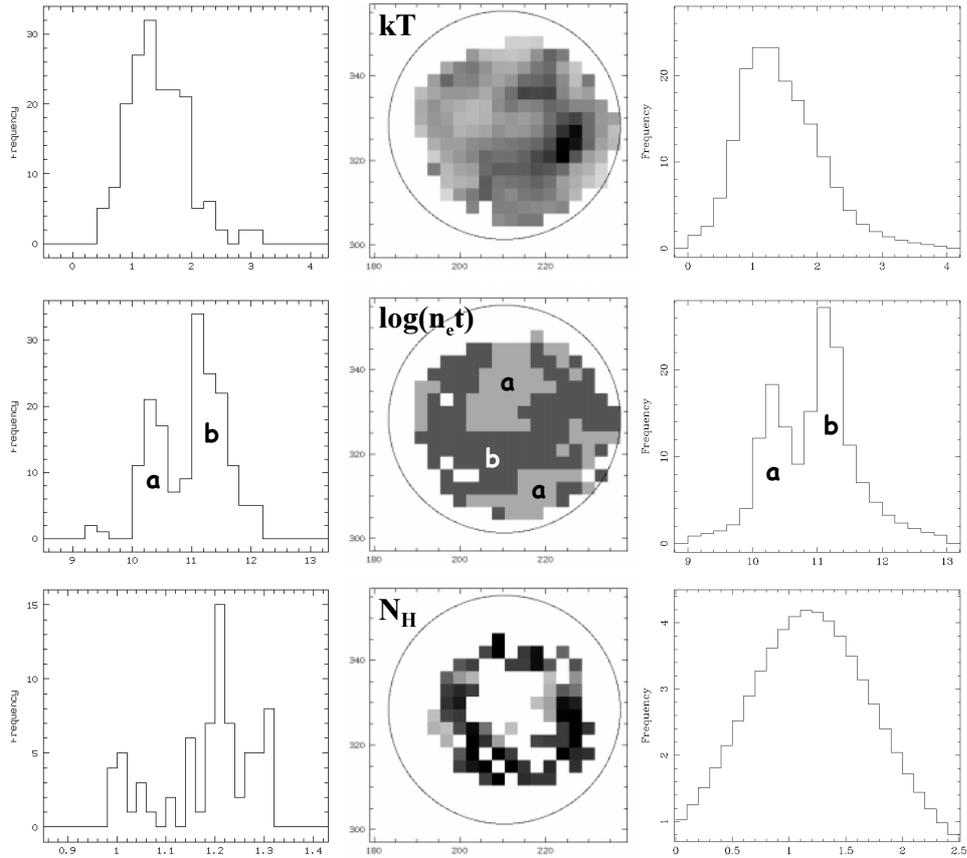
**Fig. 3.** Spatially resolved spectra of the pixels marked with a letter in Fig. 2. For each frame, top panel refers to data and folded model, while bottom panel refers to the fitting residuals in unit of data-to-folded model ratio

thermal models with different physical and chemical parameters, interpreted as an indication of the presence of two different components in the remnant. The only exception is given by the analysis of *Tenma* data (Tsunemi et al. 1986). In particular, *BeppoSAX* unresolved spectrum has been modeled with two NEI components respectively at the temperature of 1.25 and 3.8 keV. Our spatially resolved spectroscopic analysis allowed us to get statistically significant fits using only one thermal component per pixel with temperatures compatible with the one relative to the

cooler component in Favata et al. (1997). This means that if two thermal components are indeed present, they are spatially separated.

### 5.1. $kT$

$kT$  values are distributed in the range 0.9–3.2 keV, with most of the values in the rather narrow 1–2 keV range. The distribution broadens slightly if we take into account the uncertainties on the derived values (Fig. 4, right panel).



**Fig. 4.** NEI plasma model: frequency and spatial distribution of  $kT$ ,  $\log(n_e t)$ , and  $N_H$ . The superimposed circle is only for positional reference. The right panels show the frequency distribution of the same data with errors taken into account (see text)

The shape of the continuum does not vary considerably across the remnant. We notice three exceptions: (i) a well defined minimum NE of centre with  $kT \sim 0.9$  keV, (ii) a maximum W of centre with  $kT \sim 3$  keV, correlated with the relative maximum in total emission indicated as “C” in Fig. 2, and (iii) low values at the edge of the remnant (with the exception of the southern edge). A hot western component was also found by Vink et al. (1996) analyzing their ASCA data of Cas A.

It is well known from the physics of strong shocks through a gas with adiabatic index  $\gamma = 5/3$ , that electron temperatures in the range 1–2 keV can be reached by post-shocked matter when the shock velocity is of the order of 900–1300 km s<sup>-1</sup>. From observations in the optical (Reed et al. 1995) and in the X-rays (Holt et al. 1994; Vink et al. 1998) we can infer that the main shock of Cas A is presumably running now more than 2 times faster, producing post-shock temperatures over 4 times higher than observed. Even in the hypothesis of a recent deceleration due to the interaction with a denser ISM, we should see in the  $kT$  map the signature of such a deceleration in the form of a substantial increase of  $kT$  at the edge(s) of the remnant. On the other hand, higher resolution X-ray images clearly show a strong surface brightness contrast between the main shock and the bright ring that dominates the X-ray and radio emission. Nowhere in our 182 locations across the remnant we detect high temperatures,

strongly indicating that the main shock is now interacting with a very tenuous ISM, generating an emission that is too faint to be detected in our observations.

Although the morphology of the distribution of  $kT$  values is certainly of some importance, the real question to ask is: why do we detect a  $kT$  in the range 1–2 keV? A correlated question is: why do we see matter enriched in heavy elements (Si through Ni) emitting in the X-rays at such temperatures? A similar question will be asked further on in a different context and for different reasons. What we know about the hydrodynamics of the explosion inside a star is rather well established: the shells of the exploding star closer to the mass-cut (e.g. the mass coordinate separating what will be ejected from the stellar remnant), carrying out the yield of explosive nucleosynthesis, reach extremely high temperatures fractions of a second after the beginning of the explosion inside the star. Few hours later, the main shock is still inside the star, at the base of the photosphere, ready to break through giving rise to the standard SuperNova event in the  $V$  and  $UV$  bands. At the same moment, those inner shells pulled out by the main shock have expanded almost to the surface of the star. But the key point is that they have already cooled down to temperatures lower than 0.01 keV.

If emission is dominated by the ejecta (but Borkowski et al. 1996, suggest a different interpretation), how to re-heat those ejected shells? The simplest answer can be

found in Astrophysics textbooks and reviews (e.g. McKee & Truelove 1995): ejected matter acts like a piston, a secondary shock forms, running back (in lagrangian coordinates) into the ejecta, re-heating it to temperatures lower than those produced by the main forward shock. We just note *en passant* that nobody thought of shifting the onset of secondary shocks to places where the probability of formation is certainly high (inside the star), and to the right time (hours before the SuperNova event).

Being this not the right place to question the applicability of this model to this particular case, we limit ourselves to suggest a different, more straightforward scenario, based on what we observe also in other wavelength ranges. The basic ingredients of our scenario are azimuthal secondary shocks to convert a small fraction of the bulk kinetic energy ( $\sim 10^{51}$  erg) of “polluted” shells into thermal energy. Only few % of kinetic energy need to be converted by azimuthal shocks into thermal energy to set up the scene for what we will observe  $10^{10}$  s later. Meanwhile, vorticity will help forming blobs or knots of frozen local composition and scrambling their relative positions.

The scenario is the following: at time  $t \simeq 0$  (plus or minus hours, does not really matter) we deal with shells of material enriched in heavier elements that are rather cold. Energy is mostly in the form of kinetic energy. Their bulk motion is subsonic with respect to adjacent shells, but supersonic with respect to the value of the sound speed in the shells. This is a very unstable situation. Any element of matter with a random azimuthal velocity component will be shocked (or compressed) by radially moving adjacent matter. Kinetic energy will be transformed into thermal energy, and vorticity will set in. Of course the complexity of this basic scenario, based on azimuthal shocks in a cold supersonic flow, can be increased *ad libitum* adding rotation, and in general any non spherically symmetric effect.

## 5.2. $N_{\text{H}}$

Due to the limited extension towards low energies of the MECS well calibrated band ( $E \geq 1.6$  keV) we cannot determine  $N_{\text{H}}$  with reasonable accuracy. Taking into account the uncertainties in the determination of  $N_{\text{H}}$  (see Fig. 4, right panel) we can say that our results are consistent with those of Kehoane et al. (1996). Moreover, we fitted one of the most significant spectra (namely, the spectrum marked with “A” in Fig. 3, corresponding to the pixel  $X = 203$ ,  $Y = 321$  in Table 4<sup>5</sup>) fixing  $N_{\text{H}}$  to three different values; the results, as listed in Table 2, show that no significant variations in any of the  $N_{\text{H}}$  sensitive spectral parameters are detected.

## 5.3. $n_e t$

The ionisation parameter  $n_e t$  controls the non equilibrium ionisation status of the emitting plasma. Because we

**Table 2.** Test for the spectrum marked with “A” in Fig. 3 of the sensitivity of the derived parameters to  $N_{\text{H}}$ . Quoted errors are one standard deviation

$N_{\text{H}}$ ( $10^{22}$ cm $^{-2}$ )	0.7	1.3	1.9
$kT$ (keV)	$1.6 \pm 0.26$	$1.46 \pm 0.29$	$1.4 \pm 0.22$
$\log(n_e t)$ (cm $^{-3}$ s)	$11.25 \pm 0.2$	$11.3 \pm 0.22$	$11.3 \pm 0.2$
[Si]	$5.9 \pm 1.35$	$6.46 \pm 1.72$	$8 \pm 2$

express  $n_e t$  in units of cm $^{-3}$  s and its value can vary by orders of magnitude, in Table 4, in Fig. 4 and in this discussion we will always refer to values of  $\log(n_e t)$ . For the elements considered in this paper (Si through Ni) a low value of the ionisation parameter (e.g.  $\log(n_e t) < 10.5$ ) indicates an ionisation structure far from equilibrium, while the structure is practically in equilibrium for values greater than 12. In the fitting procedure  $\log(n_e t)$  is allowed to vary in the range 6–12.5.

The frequency distribution of  $\log(n_e t)$  values is clearly bimodal, with lower boundaries at about 10 and 11. Because an upper limit for  $t$  is the age of the remnant (assumed to be 320 y, e.g.  $\log(t) = 10$  in our units), this can give us a good handle in understanding the density distribution within the remnant. The simplest hypotheses concerning the density distribution are:

case (i): we fix  $t$  to infer the distribution of  $n_e$ . We are observing matter shocked (once or more than once)  $10^{10}$  s ago, that is during the early phases of the explosion. Then the distribution of the ionisation parameter actually reflects the distribution of the electron density, shifted by 10 orders of magnitude. If this hypothesis is true, we detected a bimodal density distribution, with a first component (marked *a* in Fig. 4) with  $n_e$  in the range 1–10 cm $^{-3}$  and a second component (marked *b* in the same figure) about 10 times denser. All in all  $n_e$  (actually: the electron density averaged over our pixel volume) would cover the range 1–> 100.

case (ii): we fix the distribution of  $n_e$  to infer the distribution of  $t$  (e.g. of the main shocking episodes). The sharp boundary at  $\log(n_e t) = 11$  indicates that in the remnant is present only one component (*b*) with (pixel averaged)  $n_e$  in the range 10–100 cm $^{-3}$  and over that was shocked  $10^{10}$  s ago. From the boundary at  $\log(n_e t) = 10$  we can infer that a consistent fraction of this same matter (with  $n_e \geq 10$  cm $^{-3}$ ) was suddenly re-shocked  $10^{10}$  cm $^{-3}$  s /  $10$  cm $^{-3}$  =  $10^9$  s  $\simeq$  32 y ago (that is in the mid sixties, when Cas A was already well observed). There could have been two different causes for this sudden re-shocking of part of the remnant: external, e.g. matter at the edge of the remnant hit pre-existing clouds or condensation; internal, e.g. blobs or fingers of material inside the remnant collided and re-heated. What we observe in Fig. 4 is that this hypothetically re-shocked material is located mostly at the centre of the remnant and north of it, with few other pixels at the edges. But no particular anomalies in  $kT$  are found at the corresponding locations that could be interpreted as signatures of a very recent re-shocking.

<sup>5</sup> Table 4 is available only in electronic form at <http://www.edpsciences.org>

These first two cases are the extreme possibilities in which we fix either  $t$  or  $n_e$ .

(iii) a third intermediate possibility is that matter is distributed in components  $a$  and  $b$  as in (i), plus an additional component  $c$  of density higher than component  $b$  by a factor  $n$ , that was then re-heated  $age/n$  years ago. Factor  $n$  must be significantly greater than 1 in order to consider this case different from case (i).

The above hypotheses have an obvious impact on the determination of the total mass  $M$  of the remnant:  $M(ii)$  can be up to 4 times  $M(i)$ , and  $M(i) < M(iii) < M(ii)$ .

We favour hypothesis (i) on the basis of the following considerations:

- (1) case (i) is not in contrast with the other results of our analysis. On the other hand we don't see evidence for the re-heating implied by cases (ii) and (iii);
- (2) the observed values of  $kT$  are much too low to indicate the presence of shocked ISM (i.e. matter not ejected by the progenitor star in its pre-SN life). These values imply that what we are observing is just matter belonging to the progenitor star spread out in a volume of about 3 pc in radius;
- (3) the lowest possible total mass of the (X-ray) emitting gas computed according to hypothesis (i) is already in the range 13–30  $M_{\odot}$ , reasonable masses for the progenitor star. By the way, X-ray emitting masses in this range were already suggested by Murray et al. (1979) and Fabian et al. (1980).

Combining the two last considerations, one sees that no much room is left for hypothesis like (ii) or (iii), that will lead to total masses exceedingly large for the progenitor star.

If we then retain only hypothesis (i), component  $a$  could be associated to the CSM, and the denser component  $b$  to the disrupted star. We should then detect a correlation between the spatial distribution of component  $b$  and that of overabundances. Anticipating the results of the discussion of Fig. 5, we can say that this hypothesis is not in contradiction with the maps of abundance values shown in Fig. 5, with the only exception of the northern sector of the remnant.

Another interesting consequence concerns the ratio by mass of these two components. Although components  $a$  and  $b$  are spread over  $\sim 40\%$  and  $\sim 60\%$  of the pixels, respectively, the mass ratio  $M(a)/M(b) = M(\text{CSM})/M(\text{ejecta})$  is  $\leq 1/10$ . This ratio is much smaller than in previous estimates (see Vink et al. 1996; Borkowski et al. 1996). In other words, our results seem to suggest that the progenitor star lost in its pre-SN phase a small fraction of its mass. If Cas A was an underluminous SN event, it wasn't because it lost most of its envelope before the explosion. Other explanations should be found: a moderate initial mass or a high mass-cut at the moment of the explosion, if component  $a$  represents the total CSM mass, or, alternatively, we are detecting only a fraction of the total CSM, which is actually distributed over a much larger volume than the remnant. This would be the case if

much of the mass loss occurred when the progenitor was a blue supergiant.

We also remark that the denser component is not distributed in our map (in Fig. 4) as it should if matter had been ejected in spherical symmetry. On the contrary, the shape of the distribution of component  $b$  in Fig. 4 could support the hypothesis of a toroidal ejection with an axis of symmetry oriented from NE to SW. Markert et al. (1983) were the first to notice from X-ray spectra of Cas A, taken with the crystal spectrometer on board the *Einstein* Observatory, that the northwest part of Cas A is red-shifted with respect to the southeast, with a dividing line running then from NE to SW. The same symmetry was found by Holt et al. (1994) in their ASCA data, from which they could build a Doppler contour map of Cas A showing that the SE part of the remnant was blue-shifted and the NW region was red-shifted. The torus could be by now broken in a number of different concentrations, giving rise to the SE emission enhancement (running towards us) and other in the north and west (running away from us).

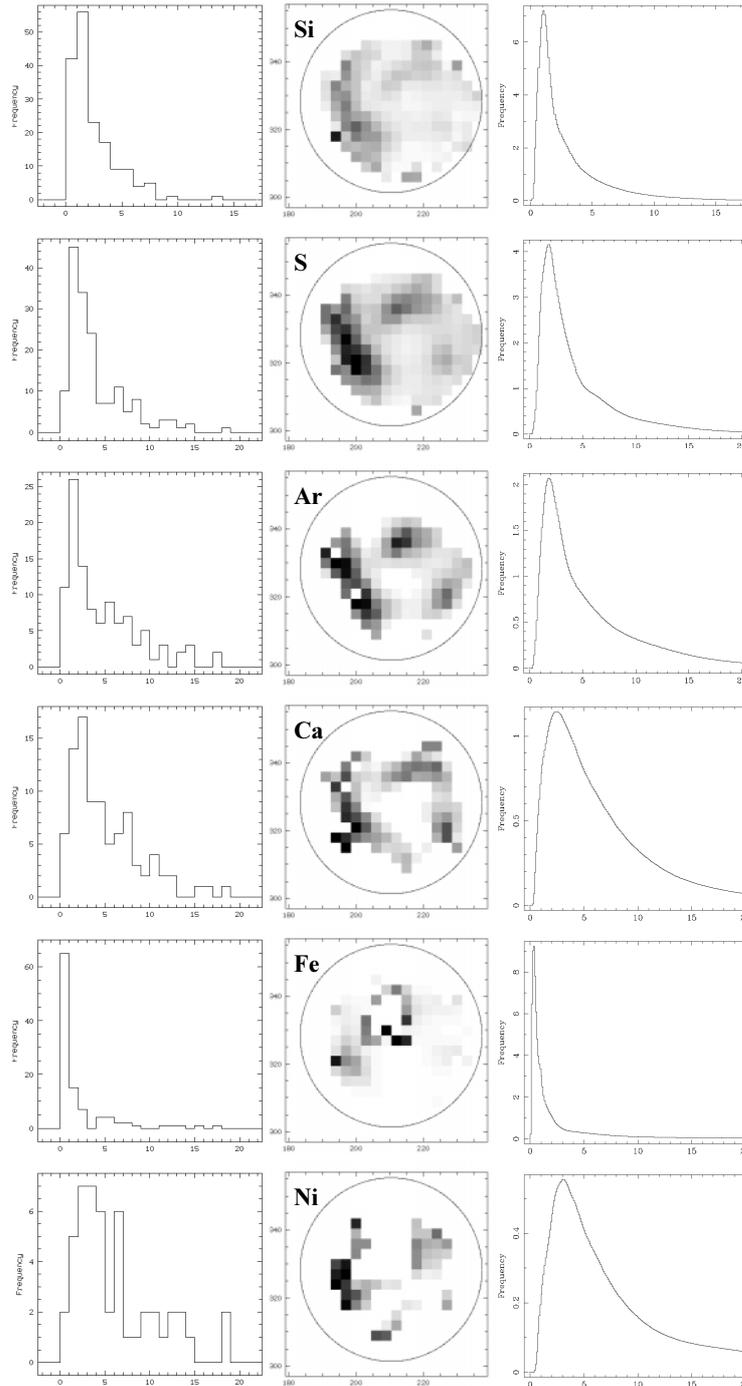
#### 5.4. Abundances

Abundances were derived, with varying degree of accuracy, for Si in 167 out of the 182 pixels, for S in 164 pixels, for Ar in 106 pixels, for Ca in 90 pixels, for Fe in 105 pixels, and for Ni in 49 pixels. The determination of Si and Ni abundances is affected by the position of the corresponding blends at the opposite extremes of our energy window. The fit of the Si blend is affected by the inability to constrain reasonably well the value of  $N_{\text{H}}$ , while the Ni blend is located in a region of the spectrum where the count rate is the lowest, although the presence of a Ni line is not related to the statistics of the spectrum: see spectra A and B in Fig. 3. The list of elements in order of decreasing global accuracy of the corresponding abundance determination is: S, Fe, Si, Ar, Ca, and Ni.

The frequency distributions of abundance determination (with respect to solar, by number) are shown in the left panels of Fig. 5. The right panels map the spatial distribution of these values, with the same intensity scale for all six panels. The same reference circle centered at the expansion centre (as defined by Reed et al. 1995) is shown in all panels for sake of comparison of the different distributions.

Now, we know that Cas A is located not much further away from the centre of the Galaxy than our Sun, and was generated by the explosion of a massive star born between about 3 to 10 My ago. In terms of the time scale of galactic evolution, this means “now”, and today's abundances of the ISM at the location of Cas A are basically solar. On all this there is a general consensus. As well as on the obvious expectation to detect overabundances of heavy elements (in our case: Si to Ni) with respect to solar values, as found in the optical knots.

But from common knowledge (certainly still limited) of explosive nucleosynthesis the following question is fully



**Fig. 5.** NEI plasma model: frequency and spatial distribution of Si, S, Ar, Ca, Fe, and Ni overabundances. The superimposed circle is only for positional reference. The right panels show the frequency distribution of the same data with errors taken into account

justified: should we really expect to observe those overabundances? The answer is: yes, for  $\alpha$ -elements, and: not necessarily, for Fe and Ni. Suppose the extreme hypothesis of a complete mixing, above the mass-cut, of all the matter belonging to the progenitor star. The production of  $\alpha$ -elements is such that, unless stellar matter is diluted in a much greater mass of ISM, we do expect to observe overabundances of Si through Ca. On the other hand, depending on the precise value of the mass-cut, overabundances of Fe and Ni could well pass undetected.

A more likely scenario, based on high spatial resolution studies of this remnant in the optical (and now also in the X-rays), is that Cas A is a remnant composed by scrambled “blobs” or knots of frozen local material more than a mixed distribution of matter. In this case we should re-formulate our answer to the previous question, saying that: yes, we do expect overabundances of heavy elements, but only in the location of parcels or blobs of ejecta keeping trace of the effects of explosive nucleosynthesis. In other locations we can only expect to find abundances

**Table 3.** Global overabundance factors by mass. Columns marked with (1) and (2) refer to different assumptions discussed in the text

Element	(1)	(2)
[Si]	3.3	3.1
[S]	5.3	5.0
[Ar]	6.3	4.1
[Ca]	6.8	4.6
[Fe]	1.5	1.4
[Ni]	7.1	3.9

that will reflect those of the ISM at the time the progenitor star was born. Actually, due to *s*-process nucleosynthesis, the Fe abundance could even be depleted below its initial value (see for instance the recent paper by Limongi et al. 2000).

From the left panels of Fig. 5 we clearly see that in the majority of the pixels we are detecting abundance values compatible with solar values. More in detail, there seems to be indications that in the location of Cas A abundances of Fe and Ni are somewhat below solar, while abundances of the  $\alpha$ -elements are above solar. As an example, for Fe we find that about 2/3 of the pixels show an average abundance  $\sim 1/2$  solar, and for S the average abundance is  $\sim 1.7$  solar in 40% of the pixels. For comparison, Borkowski et al. (1996) derived for the CSM component abundances  $0.4\times$  solar for Fe and Ni, and slightly below solar for other elements heavier than Ne.

On the other hand we do see overabundances in particular locations inside the remnant that lead to global overabundances. From the results listed in Table 4 we can compute for each element the ratio  $[X]$  between the mass fraction per unit solar mass of element *X* and the mass fraction of the same element in the sun. We can compute this overabundance by mass in two different assumptions: (1) we can consider only pixels in which we detect the element, or (2) we can consider also the contribution of all the other pixels where we assume that abundances are solar (i.e. were not altered by nucleosynthesis in the progenitor star). The global overabundance factors we find according with these two assumptions are listed in Table 3.

As we can see from the right panels of Fig. 5, all the elements are particularly overabundant in the south-east. This region is extending up to the east in Si, S, Ar, and Ca maps. At lower level of overabundance these four elements trace a sort of broken shell, more evident in the case of Ar and Ca maps; similar results were found by Holt et al. (1994) and Hwang et al. (2000).

## 6. Conclusions

We have performed the first detailed spatially resolved spectroscopy of Cas A in the X-rays, using data taken with the MECS spectrometer on board the *BeppoSAX* Observatory. The extremely well calibrated PSF in the central region of the MECS allowed us to perform a spatial

deconvolution of our data at full energy resolution. We eventually generated a data cube consisting of 182 spectra, each one referring to a squared pixel of  $24''$  in size, covering a region of  $\sim 3'$  radius around the centre of Cas A.

We performed a fit of the 182 spectra using as models a NEI plasma and a power law with fixed normalization and index to account for the high energy tail observed by *BeppoSAX*. Statistical errors on counts have been modified to take into account the systematic effects of the deconvolution procedure.

In a previous study, reasonable fits of the spatially unresolved spectrum of Cas A observed by *BeppoSAX* were obtained using two thermal models with different physical and chemical parameters, indicating the presence of two different components in the remnant. On the contrary, spatially resolved spectroscopy allowed us to get statistically significant fits using only one thermal component per pixel. This means that if two thermal components are indeed present, they are spatially separated.

We found a rather uniformly distributed  $kT$ , with a minimum in the east, a maximum in the west, and most of the values in the range 1–2 keV. No evidence is found for high  $kT$  expected from the interaction of the main shock with the ISM. No anomalies in the  $kT$  distribution are found at the edge of the remnant. We interpret this as an indication that we are detecting only matter belonging to the progenitor star.

We have then discussed how to re-heat matter to the observed values, in particular matter enriched by heavier elements, that is expected to have already cooled down to  $< 0.01$  keV at the moment of the SN event. A possibility is to invoke the standard “reverse shock” model. In this paper we have also suggested a scenario based on azimuthal secondary shocks in the supersonic flow generated by the main shock inside the star. Only a small fraction of the kinetic energy needs to be converted into thermal energy by azimuthal shocks. The peculiarity of this scenario is that azimuthal shock can be formed from hours before to years after the SN event. Actually they could still be at work now. Observations in the X-rays at much higher spatial resolution could help in clarifying the issue.

From the discussion of our results on the ionisation parameter  $n_e t$  we infer that two distinct components there exist: the first with  $n_e$  in the range  $1\text{--}10\text{ cm}^{-3}$ , the second with  $n_e$  ten times (or more) higher. The first component (called *a*) can be associated with matter ejected by the star before the SN event, while the denser component (*b*), not spherically symmetric, can be associated with matter ejected during the explosion. An important consequence of this interpretation is that the mass ratio  $M(a)/M(b)$  is  $\leq 1/10$ , indicating a progenitor star of moderate mass that lost only a small fraction of the envelope during its pre-SN life.

We also interpret the distribution of component *b* across the remnant as an indication that the explosion was not spherically symmetric, possibly toroidal.

The distribution of abundances indicates that we are detecting a CSM component with almost solar

composition, and an ejecta component enriched in heavier elements. Abundances found for  $\alpha$ -elements are consistent with the current view that Cas A was produced by the explosion of a massive star. But, as already suggested by Vink et al. (1996), low Fe overabundances can be an indication that at the moment of the explosion the mass-cut was rather high, locking most of the produced Ni<sup>56</sup> into the stellar remnant. The value of Fe abundance found in the CSM component (less than the solar one) is again in agreement with theoretical predictions.

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## Appendix A: Error treatment

Errors associated to the counts in each pixel of the final images are the combination of the canonical statistical error (squared root of the counts) and of a contribution introduced by the deconvolution that depends on the statistics, on the distribution of counts within the image, on the PSF approximation, and on the convergence criteria. We evaluated the total errors empirically assuming that:

- I - deconvolved images relative to energies closer than the instrumental spectral resolution must be identical within the errors;
- II - in each spectrum accumulated from the deconvolved images, the spread of the points respect to a model that well describe the overall behaviour (giving a flat distribution of the residuals) must be within the errors.

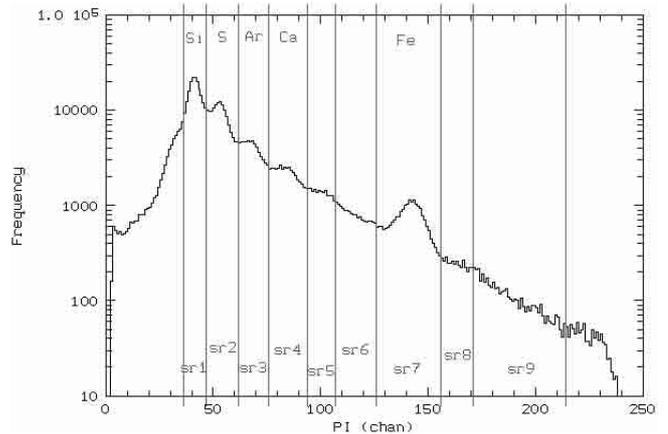
Each one of these two hypotheses can separately be used to evaluate the errors, and the results derived from one hypothesis must automatically satisfy the other one. We have implemented two procedures, one based on the images (I) and the other on the spectra (II), and then we have compared the two sets of results.

### A.1. Method I: Images based

Following the hypothesis that deconvolved images close in energy must be identical within the errors, we identified 9 energy ranges, most of which are coincident with the source lines, where the spatial distribution of the source emission does not vary with energy. The boundaries chosen for these 9 spectral regions are shown in Fig. A.1.

Taking into account only the well deconvolved spatial area, we have produced 9 templates (one for each spectral region) adding all images in a range and normalizing the source counts in the summed image. The deconvolved images of each PI in the same spatial area have then been compared to the relative template scaled for the source counts in that area by using the variable  $\chi_{\text{st\_ima}}^2$  and assigning the statistical errors to the pixel counts:

$$\chi_{\text{st\_ima}}^2 = \frac{1}{N_{\text{dof}}} \sum_{i,j} \frac{(D_{i,j} - T_{i,j})^2}{D_{i,j}}$$



**Fig. A.1.** The spectral regions used to estimate errors superimposed on data from the ME3 unit, Op2990

where  $D_{i,j}$  refers to the counts in the pixel  $(i, j)$  of the deconvolved PI image,  $T_{i,j}$  is the corresponding value in the related template, and  $N_{\text{dof}}$  (degrees of freedom) is the number of pixels on which  $\chi_{\text{st\_ima}}^2$  is evaluated. Only pixels with more than 15 counts have been included in the computation. The variable  $\chi_{\text{st\_ima}}^2$  is related to the final errors through a few simple assumptions:

- the final errors  $Err$  are related to the statistical ones by:

$$Err = \alpha_{\text{ima}} Err_{\text{st}}$$

where  $Err_{\text{st}}$  are the statistical errors and  $\alpha_{\text{ima}}$  is the correction factor, always greater than 1;

- the deconvolution introduces errors that at first order do not depend on the position in the image, i.e.  $\alpha_{\text{ima}}$  is constant over each PI image.

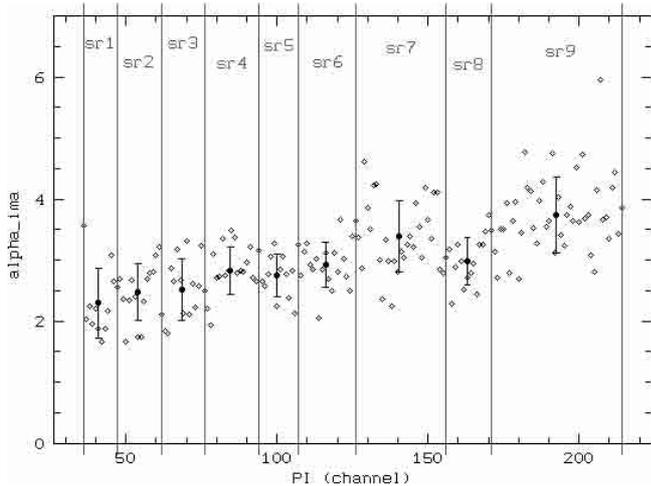
Following these two assumptions, and being the expected value of the reduced  $\chi_{\text{red}}^2$  equal to 1, we obtain:

$$\alpha_{\text{ima}} = \sqrt{\chi_{\text{st\_ima}}^2}$$

Figure A.2 plots the  $\alpha_{\text{ima}}$  values found at the different PI channels. The average value of  $\alpha_{\text{ima}}$  in each template region is also shown with its relative error (root-mean-square deviation).

### A.2. Method II: Spectra based

We implemented a second procedure following the hypothesis that in each spectrum the spread of the points respect to a model, that well describe the general behaviour, must be within the errors. We chosen to analyze only spectra having a total number of counts greater than  $4 \cdot 10^4$ : in these spectra in fact the statistics of most of the energy channels is sufficiently high to allows us to use the  $\chi^2$  test without rebinning. Actually, some of the high energy channels present very low or zero counts; however, for homogeneity and to easily automatize the procedure, we did not rebin spectral channel.



**Fig. A.2.** The  $\alpha_{\text{ima}}$  correction factor as a function of the PI channels. Each point refers to a single image. The average value of  $\alpha_{\text{ima}}$  in each template region is also shown with its relative error (root-mean-square deviation)

We fitted the selected spectra (a total of 24) using a model that, although it does not correctly describe the emission mechanism of the source, gives a flat distribution of the residuals for all the spectra. The fitting model is composed by a power-law plus 9 lines. All parameters are left free to vary but the line widths that are kept fixed to zero. The MECS response matrix and the background spectrum are the same used for the spectral analysis described in Sect. 4. The spread of the residuals with respect to the zero level is not uniform over the whole range: high energy channels present a wider spread respect to the low energy ones.

To compute the final errors we need to set some assumptions:

- the final errors  $Err$  are related to the statistical ones by the same formula as in Method I (Sect. A.1):

$$Err = \alpha_{\text{spec}} Err_{\text{st}}$$

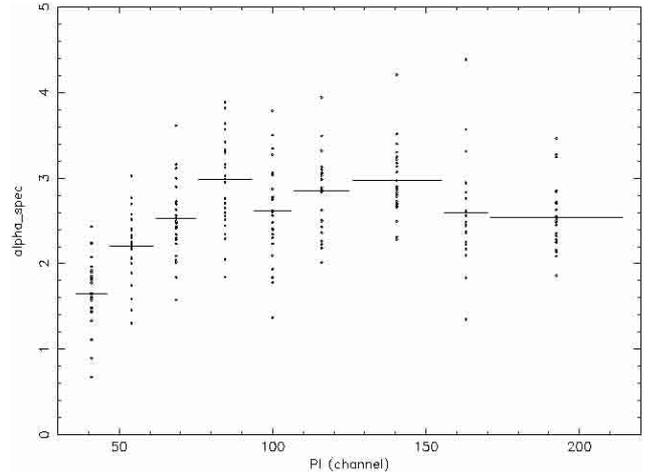
where  $Err_{\text{st}}$  are the statistical errors and  $\alpha_{\text{spec}}$  is the correction factor;

- the value of  $\alpha_{\text{spec}}$  is a function of the PI channel and does not depend on the counts detected in it;
- $\alpha_{\text{spec}}$  can be approximated to a step function of PI being constant within each given energy range.

Note that these assumptions are equivalent to those used in the image based method (Sect. A.1) and are based on the hypothesis that  $\alpha_{\text{spec}}$  mainly depends on the statistics of the deconvolved images.

Following these assumptions we divided our spectra in 9 energy ranges coincident with the ones used in Method I (see Fig. A.1). Within each of these ranges  $\alpha_{\text{spec}}$  is assumed constant and related to  $\chi_{\text{st\_spec}}^2$ , computed using the statistical errors, by the following formulas:

$$\chi_{\text{st\_spec}}^2(sr_k) = \frac{1}{N_{\text{dof}}(sr_k)} \sum_{PI} \frac{(F(PI) - C(PI))^2}{C(PI)}$$



**Fig. A.3.** The  $\alpha_{\text{spec}}(sr_k)$  correction factor in the 9 defined energy bands. Horizontal bars denote their mean values. Each point in any  $sr_k$  band refers to one of the 24 selected spectra

$$\chi_{\text{red}}^2(sr_k) = \frac{1}{\alpha_{\text{spec}}^2(sr_k)} \chi_{\text{st\_spec}}^2(sr_k).$$

By assuming  $\chi_{\text{red}}^2(sr_k) = 1$ , we obtain

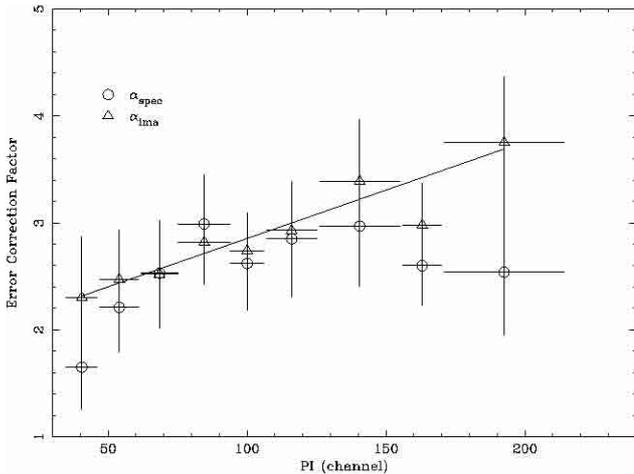
$$\alpha_{\text{spec}}(sr_k) = \sqrt{\chi_{\text{st\_spec}}^2(sr_k)}$$

where  $PI$  is ranging within the given  $sr_k$  energy range,  $F(PI)$  are the counts relative to the fitting model in the PI channel, and  $C(PI)$  are the counts detected in the same channel. Note that  $C(PI)$  corresponds to the value of  $D_{i,j}$  (used in Method I) at the given PI channel, being  $i, j$  the pixel where the spectrum has been accumulated.

We evaluated the contribution to the  $\chi_{\text{red}}^2$  in the 9 defined energy bands for each selected spectrum. Following the assumption that  $\alpha_{\text{spec}}$  is only function of the PI, we averaged the 24 values obtained in each of the  $sr_k$  regions assuming as error the root-mean-square deviation from the mean. Figure A.3 shows the computed  $\alpha_{\text{spec}}(sr_k)$  together with their mean value averaged over the 24 analyzed spectra.

### A.3. Correction factor

Figure A.4 shows the comparison between the correction factors  $\alpha_{\text{ima}}$  and  $\alpha_{\text{spec}}$  evaluated through the two methods, image and spectra based, respectively. The values of  $\alpha_{\text{spec}}$  and  $\alpha_{\text{ima}}$  are always compatible within the errors. The higher discrepancy can be observed in the first ( $sr_1$ ) and last ( $sr_9$ ) point. In the  $sr_1$  interval,  $\alpha_{\text{ima}}$  is higher than the corresponding  $\alpha_{\text{spec}}$ ; a possible explanation is that this range contains the most statistical images and the hypothesis that the correction factor is independent from the counts in the pixel could not be sufficient: the  $\chi^2$  averaged over the whole image is higher than the one relative to the most statistical pixels. In the case of the  $sr_9$  interval,  $\alpha_{\text{spec}}$  is lower than the corresponding  $\alpha_{\text{ima}}$  being strongly affected by the PI channels with low or zero counts (these are not included in the computation of  $\alpha_{\text{ima}}$ ).



**Fig. A.4.** The  $\alpha_{\text{ima}}$  and  $\alpha_{\text{spec}}$  correction factors at different PI channels. Error bars correspond to the root-mean-square deviations. The solid line represents the model used to derive the error correction factor, fitting the  $\alpha_{\text{ima}}$  points

We then computed the error correction factor from the relation  $\alpha_{\text{ima}}$  vs. PI channel number (Fig. A.2). Fitting these points with a line we get the following parameters:

$$\alpha_{\text{ima}} = 1.95 + 9.0447 \cdot 10^{-3} \times PI. \quad (\text{A.1})$$

The rms of the points respect to the line is 0.5.

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