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Native vacancies in Te-doped (5×10^{16}–5×10^{18} cm^{-3}) GaAs were investigated by means of positron lifetime and Doppler-broadening coincidence spectroscopy. The experimental data were related to theoretical calculations of the positron lifetime and the annihilation momentum distribution. Monovacancies were observed in all Te-doped GaAs samples under study. It will be shown that they can directly be identified to be Ga-vacancy–Te_{As}, donor complexes. These complexes are the dominating type of vacancy defects in the doping range under observation. [S0163-1829(99)07327-0]

Vacancies may determine important properties of semiconductor materials like GaAs. They mediate, e.g., dopant diffusion or reduce the density of free carriers. The detailed microscopic identification of vacancies and vacancy complexes in GaAs was, however, found to be difficult. Most theoretical calculations as well as diffusion studies indicate a dominating role of negative Ga vacancies in n-doped GaAs. In contrast, a recent calculation showed that the As vacancy could be an abundant defect in highly n-doped GaAs due to a low formation energy. Moreover, pairing of acceptorlike vacancies with positive donors is expected due to Coulomb attraction. Evidence for such complexes is based on photoluminescence, infrared absorption, and theoretical considerations of the doping behavior. Recently, scanning tunneling microscopy (STM) directly identified Si_{Ga}-donor–Ga-vacancy complexes on cleavage planes of highly Si-doped GaAs (Ref. 8) later shown to be present with the same density in the bulk. To our knowledge no such direct identification has been obtained so far for other n-dopants, e.g., tellurium. Positron annihilation, however, is directly sensitive to vacancies. Positron lifetime spectroscopy indeed showed the existence of native vacancies in n-doped GaAs. However, the positron lifetime measurement (which probes mainly the open volume) is not alone able to identify the defects as a given isolated arsenic or gallium vacancy, or as a vacancy-impurity complex.

One possibility to overcome the difficulty mentioned above is the investigation of the positron annihilation momentum distribution. The high momentum part of this distribution can be used to identify the chemical surrounding of the annihilation site. This is based on the fact that tightly bound core electrons with high momenta retain their element-specific properties even in a solid. This allows the identification of vacancies and vacancy-impurity complexes, especially when measurements are correlated to calculations of the momentum distribution. The coincident detection of both 511-keV γ quanta from single annihilation events allows the observation of the high momentum annihilation distribution due to a strong reduction of the disturbing background. Ga vacancies in highly Si-doped GaAs were identified using that Doppler-broadening coincidence technique. The experiment could, however, not decide whether the vacancies are isolated or a part of a complex because the Si_{Ga} donor on the second nearest site is not expected to contribute much to the annihilation. Thus, the identification of dopant-vacancy complexes in GaAs by positron annihilation is still an open question.

Tellurium is incorporated in the As sublattice only. If pairing with neighboring Ga vacancies occurs, a measurable contribution to the annihilation is expected, i.e., an identification of this complex could be possible. However, the momentum distribution for the vacancy cannot be undoubtedly determined when the fraction of trapped positrons (η) is unknown. Therefore we use correlated positron lifetime measurements to obtain η. The experimental results are compared with theoretical calculations of the annihilation characteristics to obtain a safe interpretation.

The samples studied were cut from Te-doped GaAs crystals grown by the liquid encapsulated Czochralski technique. The carrier concentration was n = 5×10^{16}, 5×10^{17}, 1.5×10^{18}, and 5×10^{18} cm^{-3}. The crystals were investigated as-received as well as annealed at 1100 °C under high (5.6 bar) arsenic pressure. The annealing conditions were chosen to maintain arsenic-rich stoichiometry and result in a maximum vacancy density for a given crystal. The material with n = 5×10^{16} cm^{-3} was not annealed. Highly Si-doped GaAs ([Si] = 1×10^{19} and 4×10^{19} cm^{-3}) was investigated for comparison since V_{Ga-Si_{Ga}} complexes were identified by STM in samples from the same wafers. Positron lifetime spectroscopy was performed using a conventional system (time reso-
The ions increase of increasing temperature positrons are detrapped from the ions relaxation alone does not allow their detailed identification. With the bulk lifetime ($t_\text{b}$) in each spectrum. The average positron lifetime ($t_\text{av}$) was characterized by the line-shape parameter $W$. The intensity of the core annihilation distribution was observed by coincident spectroscopy at 300 K using two Ge-detectors in a collinear geometry. The intensity of the core annihilation was described by the line-shape parameter $W$, defined as the integrated intensity in the region (15–20) $10^2$ counts were collected for this study is that $t_\text{d}$ is temperature $T$ dependence of positron trapping are similar in all samples, e.g., the binding energy of positrons to the Rydberg states was $E_b=(65\pm20)\text{meV}$ as was found earlier, only the concentrations of the ions and vacancies relative to each other. Note that the decrease of $t_\text{av}$ with decreasing $T$ in the highest doped sample can be similarly described (only $E_b$ was fixed in the fit). Positron trapping at vacancies is practically saturated here. Thus, $t_\text{av}$ reflects the slight decrease of the defect-related lifetime, $t_\text{d}$ with temperature [Fig. 1(b)]. This might be attributed to lattice expansion too, although the effect is larger than in the reference. The important result for this study is that $t_\text{av}=254\pm3$ ps at 300 K is the same in all samples and exhibit the same temperature dependence. This points towards a similar nature of the vacancies, assigned to be monovacancies due to the lifetime of 254 ps.

We then have to answer the question whether the vacancies observed in GaAs:Te do indeed all have the same microscopic structure. This can be achieved by checking the linearity between $t_\text{av}$ and $W$ parameter. The fraction of positrons annihilating in vacancies is given by $\eta=(W_b-W)/(W_s-W_d)=(t_\text{av}-t_\text{d})/(t_\text{av}-t_\text{b})$. Thus, the measured $W$ depends linearly on $t_\text{av}$ if the defect density changes (i.e., $\eta$) and not the defect type (i.e., $t_\text{d}$ or $W_d$). The $W-t_\text{av}$ analysis is shown in Fig. 2. The data for differently doped GaAs:Te show a linear behavior (indicated by the solid line) within the errors. Thus, the vacancy type is identical in all Te-doped samples. Moreover, only one type of vacancies exists in all samples since the line runs through the bulk value ($t_\text{b}=229\text{ps}$, $W_b=1$). The presence of negative ions does not influence this result because their annihilation parameters agree with that of the bulk. The characteristic $W$ parameter for the vacancy in GaAs:Te can be determined from $t_\text{d}$.
The core annihilation is more intense in orders of magnitude, making a detailed comparison unnecessary since the momentum distribution itself spans several decades. The data are normalized by taking the ratio to a GaAs:Zn reference. This indicates a different defect type. The detailed observation of the annihilation momentum distribution shown below allows to relate the differences to the presence of a Te atom close to a Ga vacancy in GaAs:Te.

In Fig. 3(a) the high-momentum part of the annihilation momentum distribution is shown for the vacancies in GaAs:Te and GaAs:Si. The spectra (total area $3.5 \times 10^7$ counts) were brought to unity and scaled to full trapping at the vacancies before normalization. Lines result from smoothing and serve to guide the eye only. (b) Ratio of the momentum density to bulk GaAs for different vacancies in GaAs from theoretical calculations. The curves for $V_{Ga}-Te_{As}$ and $V_{Ga}-Si_{Ga}$ are highlighted to emphasize the good agreement with the respective experimental data in GaAs:Te and GaAs:Si. The theoretical curves are not accurate for $p_L < 15 \times 10^{-3} \text{m}_0 \text{c}$ (Ref. 14) and hence are omitted.

The theoretical curves are not accurate for $p_L < 15 \times 10^{-3} \text{m}_0 \text{c}$ (Ref. 14) and hence are omitted. The ratio of the momentum density to bulk GaAs normalized by taking the ratio to a GaAs:Zn reference for the vacancies in GaAs:Te and GaAs:Si. The theoretical curves are not accurate for $p_L < 15 \times 10^{-3} \text{m}_0 \text{c}$ (Ref. 14) and hence are omitted.

The core annihilation is more intense at $V_{As}$ than at $V_{Ga}$ (Ref. 14). The As-3$d$ electrons ($Z = 33$) are more tightly bound, i.e., the momentum distribution is broader and the intensity of the core annihilation is reduced. Positron annihilation at the $Si_{Ga}$-$V_{Ga}$ complexes in GaAs:Si occurs mainly with $3d$ electrons from As. Thus, the momentum distribution should be broader compared to the bulk. This is, in fact, observed in Fig. 3(a). In contrast, at As vacancies the momentum distribution should be narrower and more intense because annihilation occurs mainly with the Ga-3$d$ electrons.

In tellurium the main contribution to the core annihilation comes from $4d$ electrons that are less strongly bound than the As-3$d$ electrons in GaAs. They contribute therefore to the core annihilation more at lower momenta and have a steeper momentum distribution. A similar difference has been noted earlier by comparing results from bulk InP, GaSb, and GaAs (Ref. 14) and for Zn-impurity P-vacancy complexes in InP. The shape of the momentum distribution measured in GaAs:Te therefore indicates that the vacancies are neighbored by Te atoms. Because Te resides on the As sublattice, the vacancy must be on the Ga sublattice. Thus we identify the vacancies in GaAs:Te to be Ga-vacancy–Te$_{As}$ complexes. This assignment explains also the reduced positron lifetime at the vacancies in GaAs:Te since the large Te atom is expected to decrease the open volume of the neighboring Ga vacancy compared to that in the $V_{Ga}$-$Si_{Ga}$ complex.

To support our experimental findings, theoretical calculations of the annihilation characteristics were performed with the method introduced in Refs. 13 and 14. The momentum distribution and positron lifetime in GaAs in reasonable agreement with the experiment. The momentum distribution is calculated within the independent particle model for each core electron state using free atomic wave functions. The final distribution is obtained by summing up the contributions from each state weighted by the partial annihilation rates calculated within the generalized gradient approximation (GGA) of positron annihilation. The GGA scheme is also used to calculate the positron lifetimes. All calculations were performed without taking lattice relaxation into account. The theoretically calculated momentum distributions for different defects and defect complexes in GaAs (normalized to the bulk distribution) are shown in Fig. 3(b). Calculated positron lifetimes and $W$ parameters are given in Table I.

The lifetime calculated for the $V_{Ga}$-$Si_{Ga}$ complex (267 ps) is slightly longer than experimentally observed (262 ps). However, the discrepancy is rather small and reflects the similar difference between the measured and calculated bulk lifetimes of 229 and 232 ps, respectively. Moreover, the shape of the calculated momentum distribution and hence the $W$ parameter are in good agreement with the experimental results. Since the $V_{Ga}$-$Si_{Ga}$ complexes were undoubtedly identified we can use the respective data as a reference state, focusing on relative changes in the following.

The core annihilation is more intense at $V_{As}$ than at $V_{Ga}$ according to the calculations [Fig. 3(b)], reflected in the high calculated $W$ parameter $W(V_{As}) = 0.91$ (Table I). This is expected from the qualitative arguments above. It should, however, be noted that the calculation overestimates the intensity of the annihilation with Ga-3$d$ electrons, discussed in more detail below.

FIG. 3. (a) High momentum part of the positron annihilation momentum distribution (normalized by taking the ratio to a GaAs:Zn reference) for the vacancies in GaAs:Te and GaAs:Si. The spectra (total area $3.5 \times 10^7$ counts) were brought to unity and scaled to full trapping at the vacancies before normalization. Lines result from smoothing and serve to guide the eye only. (b) Ratio of the momentum density to bulk GaAs for different vacancies in GaAs from theoretical calculations.
TABLE I. Theoretically calculated positron lifetime and $W$ parameter (relative to the bulk) for different vacancies and vacancy complexes in GaAs.

<table>
<thead>
<tr>
<th>Vacancy</th>
<th>$\tau$ (ps)</th>
<th>$W_{\text{ref}}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>GaAs bulk</td>
<td>232</td>
<td>1</td>
</tr>
<tr>
<td>$V_{\text{As}}$</td>
<td>266</td>
<td>0.92</td>
</tr>
<tr>
<td>$V_{\text{Ga}}$</td>
<td>267</td>
<td>0.74</td>
</tr>
<tr>
<td>$V_{\text{Ga}}$-$\text{Si}_{\text{Ga}}$</td>
<td>267</td>
<td>0.72</td>
</tr>
<tr>
<td>$V_{\text{Ga}}$-$\text{Te}_{\text{As}}$</td>
<td>267</td>
<td>0.75</td>
</tr>
<tr>
<td>$V_{\text{Ga}}$-$\text{Ga}_{\text{As}}$</td>
<td>267</td>
<td>0.80</td>
</tr>
<tr>
<td>$V_{\text{As}}$-$\text{Te}_{\text{As}}$</td>
<td>265</td>
<td>0.91</td>
</tr>
<tr>
<td>$V_{\text{As}}$-$V_{\text{Ga}}$</td>
<td>332</td>
<td>0.62</td>
</tr>
<tr>
<td>$V_{\text{Ga}}$-$\text{Te}<em>{\text{As}}$-$V</em>{\text{Ga}}$</td>
<td>272</td>
<td>0.67</td>
</tr>
<tr>
<td>$V_{\text{As}}$-$V_{\text{Ga}}$-$\text{Te}_{\text{As}}$</td>
<td>328</td>
<td>0.62</td>
</tr>
</tbody>
</table>

detail in Ref. 14. Therefore, a slightly too high $W$ parameter may result for the As vacancy. However, the calculated positron lifetime at $V_{\text{As}}$ (which is less influenced by the difficulties of the calculation mentioned above) is similar to that of $V_{\text{Ga}}$-$\text{Si}_{\text{Ga}}$ whereas $\tau_d$ measured in GaAs:Te is lower. The calculations give thus no support to identify the vacancy in GaAs:Te to be $V_{\text{As}}$. Similar arguments apply for the (hypothe
tical) $\text{Te}_{\text{As}}$-$V_{\text{As}}$ complex which could be formed if As vacancies were present. Divacancies or divacancy-$\text{Te}_{\text{As}}$ complexes, on the other hand, have higher calculated positron lifetimes than $V_{\text{Ga}}$-$\text{Si}_{\text{Ga}}$ (Table I).

The calculated momentum distribution of isolated Ga vacancies has the same shape as the $V_{\text{Ga}}$-$\text{Si}_{\text{Ga}}$ complexes [Fig. 3(b)] only with slightly increased intensity. Moreover, the calculated positron lifetime of $V_{\text{Ga}}$ agrees with that of this complex ($\tau_d \sim 267$ ps). For these reasons we exclude isolated Ga vacancies as the ones detected in GaAs:Te. It is then interesting to look for the $V_{\text{Ga}}$-$\text{Ga}$-antisite complex which could, according to theory, be an abundant defect in Ga-rich, $n$-type GaAs. However, the shape of annihilation momentum distribution as well as the lifetime (267 ps) calculated for this defect disagrees with the experiment [Figs. 3(a) and 3(b)].

The last defect to be discussed is the $V_{\text{Ga}}$-$\text{Te}_{\text{As}}$ complex. The calculated lifetime for this complex (261 ps) is 6 ps lower, whereas the $W$ parameter (0.75) is higher than that of $V_{\text{Ga}}$-$\text{Si}_{\text{Ga}}$(0.72). This is in good agreement to the experimental results. Moreover, the momentum distribution calculated for the $V_{\text{Ga}}$-$\text{Te}_{\text{As}}$ complex is very similar to the distribution measured in Te-doped GaAs: the intensity is increased in the momentum range $p_L = (10-22.5) \times 10^{-3}$ $m_0c$ and a cross-over with the momentum distribution for $V_{\text{Ga}}$-$\text{Si}_{\text{Ga}}$ occurs at about $23 m_0c$. Thus, the theoretical calculations strongly support the identification of the vacancies in GaAs:Te to be Ga-vacancy–$\text{Te}_{\text{As}}$-donor complexes.

Finally, we address the correlation between doping and vacancy concentration suggested by this identification. The vacancy density can be most reliably estimated from $\tau_{av}$ at high temperatures where the influence of negative ions is negligible. Using the relation $[V_{\text{Ga}}$-$\text{Te}_{\text{As}}] = (\tau_{av} - \tau_0)\mu_{\text{v}}\tau_0/(\mu_{\text{v}}\tau_0)(\tau_{av} - \tau_0)$ and a trapping coefficient $\mu_{\text{v}} = 10^{15}/(77300 \text{ K})^{-1/2} S^{-1}$ (Refs. 9, 12, and 19), we obtain increasing vacancy densities of 0.2–0.4; 0.8–1; 3–4; and 10–20$\times 10^{17}$ cm$^{-3}$ with increasing doping concentration. The last value is a lower limit estimation due to saturated positron trapping. With the exception of the lowest doped sample, the ratio between vacancy and electron concentration is almost about $1/\tau_{av}$. The same relation between doping and vacancy concentration in GaAs:Te has been found earlier. In that work, the carrier compensation commonly observed in GaAs:Te (Ref. 3) was interpreted as being to a large part due to dopant-vacancy complexes. The present work confirms this interpretation by the direct identification of $V_{\text{Ga}}$-$\text{Te}_{\text{As}}$ complexes.

In summary, we applied positron lifetime and Doppler-broadening coincidence spectroscopy to study vacancies in Te-doped GaAs. We showed that the native vacancies in GaAs:Te can directly be identified to be Ga-vacancy–$\text{Te}_{\text{As}}$-donor complexes. This assignment is strongly supported by theoretical calculations. No other type of vacancies could be detected for carrier densities between $5 \times 10^{16}$ and $5 \times 10^{18}$ cm$^{-3}$, i.e., dopant-vacancy complexes rather than isolated vacancies are the dominating ones in sufficiently high $n$-doped, As-rich GaAs. It was demonstrated that even small differences in the positron signals can be used to obtain a direct and unambiguous identification of vacancies or vacancy complexes in GaAs thus helping to resolve former conflicting interpretations.

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