

# Pure Spin Current Injection in Hydrogenated Graphene Structures

Reinaldo Zapata-Peña<sup>1</sup>, Bernardo S. Mendoza<sup>1</sup>, Anatoli I. Shkrebtii<sup>2</sup>

<sup>1</sup>*Centro de Investigaciones en Óptica, León, Guanajuato 37150, México and*

<sup>2</sup>*University of Ontario, Institute of Technology, Oshawa, ON, L1H 7L7, Canada*

(Dated: April 25, 2017)

Lorem ipsum dolor sit amet, consectetur adipiscing elit. Etiam lobortis facilisis sem. Nullam nec mi et neque pharetra sollicitudin. Praesent imperdiet mi nec ante. Donec ullamcorper, felis non sodales commodo, lectus velit ultrices augue, a dignissim nibh lectus placerat pede. Vivamus nunc nunc, molestie ut, ultricies vel, semper in, velit. Ut porttitor. Praesent in sapien. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Duis fringilla tristique neque. Sed interdum libero ut metus. Pellentesque placerat. Nam rutrum augue a leo. Morbi sed elit sit amet ante lobortis sollicitudin. Praesent blandit blandit mauris. Praesent lectus tellus, aliquet aliquam, luctus a, egestas a, turpis. Mauris lacinia lorem sit amet ipsum. Nunc quis urna dictum turpis accumsan semper.

## I. INTRODUCTION

Lorem ipsum dolor sit amet, consectetur adipiscing elit. Etiam lobortis facilisis sem. Nullam nec mi et neque pharetra sollicitudin. Praesent imperdiet mi nec ante. Donec ullamcorper, felis non sodales commodo, lectus velit ultrices augue, a dignissim nibh lectus placerat pede. Vivamus nunc nunc, molestie ut, ultricies vel, semper in, velit. Ut porttitor. Praesent in sapien. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Duis fringilla tristique neque. Sed interdum libero ut metus. Pellentesque placerat. Nam rutrum augue a leo. Morbi sed elit sit amet ante lobortis sollicitudin. Praesent blandit blandit mauris. Praesent lectus tellus, aliquet aliquam, luctus a, egestas a, turpis. Mauris lacinia lorem sit amet ipsum. Nunc quis urna dictum turpis accumsan semper. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Etiam lobortis facilisis sem. Nullam nec mi et neque pharetra sollicitudin. Praesent imperdiet mi nec ante. Donec ullamcorper, felis non sodales commodo, lectus velit ultrices augue, a dignissim nibh lectus placerat pede. Vivamus nunc nunc, molestie ut, ultricies vel, semper in, velit. Ut porttitor. Praesent in sapien. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Duis fringilla tristique neque. Sed interdum libero ut metus. Pellentesque placerat. Nam rutrum augue a leo.

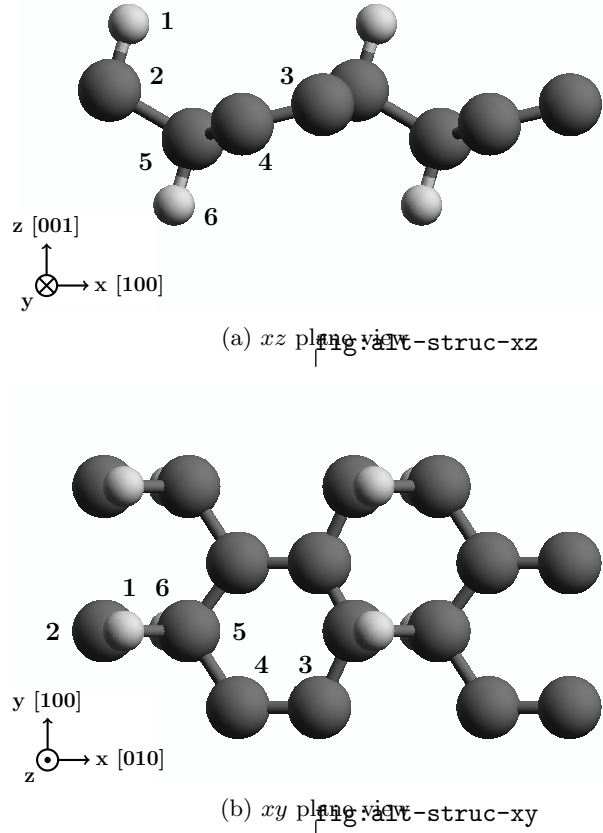


FIG. 1. Alt structure`fig:alt-struc`

Morbi sed elit sit amet ante lobortis sollicitudin. Praesent blandit blandit mauris. Praesent lec-



FIG. 2. Up structure `fig:up-struc`

tus tellus, aliquet aliquam, luctus a, egestas a, turpis. Mauris lacinia lorem sit amet ipsum. Nunc quis urna dictum turpis accumsan semper. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Etiam lobortis facilisis sem. Nullam nec mi et neque pharetra sollicitudin. Praesent imperdiet mi nec ante. Donec ullamcorper, felis

non sodales commodo, lectus velit ultrices augue, a dignissim nibh lectus placerat pede. Vivamus nunc nunc, molestie ut, ultricies vel, semper in, velit. Ut porttitor. Praesent in sapien. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Duis fringilla tristique neque. Sed interdum libero ut metus. Pellentesque placerat. Nam rutrum augue a leo. Morbi sed elit sit amet ante lobortis sollicitudin. Praesent blandit blandit mauris. Praesent lectus tellus, aliquet aliquam, luctus a, egestas a, turpis. Mauris lacinia lorem sit amet ipsum. Nunc quis urna dictum turpis accumsan semper.

Lorem ipsum dolor sit amet, consectetur adipiscing elit. Etiam lobortis facilisis sem. Nullam nec mi et neque pharetra sollicitudin. Praesent imperdiet mi nec ante. Donec ullamcorper, felis non sodales commodo, lectus velit ultrices augue, a dignissim nibh lectus placerat pede. Vivamus nunc nunc, molestie ut, ultricies vel, semper in, velit. Ut porttitor. Praesent in sapien. Lorem ipsum dolor sit amet, consectetur adipiscing elit. Duis fringilla tristique neque. Sed interdum libero ut metus. Pellentesque placerat. Nam rutrum augue a leo. Morbi sed elit sit amet ante lobortis sollicitudin. Praesent blandit blandit mauris. Praesent lectus tellus, aliquet aliquam, luctus a, egestas a, turpis. Mauris lacinia lorem sit amet ipsum. Nunc quis urna dictum turpis accumsan semper.

## II. THEORY

`sec:theory`

The equation for  $\mathcal{V}^{\text{ab}}$  for normal incidence in the *xy* plane with a polarization angle  $\alpha$  is given by

$$\begin{aligned} \mathcal{V}^{\text{ab}}(\omega) &= \frac{2}{\hbar} \frac{\mu^{\text{abxx}}(\omega)E^2(\omega)\cos^2(\alpha) + \mu^{\text{abyy}}(\omega)E^2(\omega)\sin^2(\alpha) + 2\mu^{\text{abxy}}(\omega)E^2(\omega)\cos(\alpha)\sin(\alpha)}{\xi^{\text{xx}}(\omega)E^2(\omega)\cos^2(\alpha) + \xi^{\text{yy}}(\omega)E^2(\omega)\sin^2(\alpha)}, \\ &= \frac{2}{\hbar} \frac{\mu^{\text{abxx}}(\omega)\cos^2(\alpha) + \mu^{\text{abyy}}(\omega)\sin^2(\alpha) + \mu^{\text{abxy}}(\omega)\sin(2\alpha)}{\xi^{\text{xx}}(\omega)\cos^2(\alpha) + \xi^{\text{yy}}(\omega)\sin^2(\alpha)}. \end{aligned} \quad \text{eq:vab} \quad (1)$$

For an angle  $\alpha = \frac{\pi}{4}$  this expression can be reduced to

$$\mathcal{V}^{\text{ab}}(\omega) = \frac{2}{\hbar} \frac{\mu^{\text{abxx}}(\omega) + \mu^{\text{abyy}}(\omega) + 2\mu^{\text{abxy}}(\omega)}{\xi^{\text{xx}}(\omega) + \xi^{\text{yy}}(\omega)}. \quad \text{eq:vab-90deg} \quad (2)$$

### A. Fixing velocity.

sec:theory-fixvel

Considering that we have 2D structures we fixed the velocity in the  $xy$  plane along  $x$  and  $y$  directions and we define  $|\mathcal{V}^a|$  as

$$|\mathcal{V}^a| = \sqrt{(\mathcal{V}^{ax})^2 + (\mathcal{V}^{ay})^2 + (\mathcal{V}^{az})^2}, \quad \text{eq:va-mag} \quad (3)$$

and the corresponding polar and azimuthal angles  $\theta$  and  $\varphi$  as

$$\theta = \cos^{-1} \left( \frac{\mathcal{V}^{az}}{|\mathcal{V}^a|} \right), \quad 0 \leq \theta \leq \pi, \quad \text{eq:polar-ang} \quad (4)$$

$$\varphi = \tan^{-1} \left( \frac{\mathcal{V}^{ay}}{\mathcal{V}^{ax}} \right), \quad \text{eq:azimuthal-ang} \quad 0 \leq \varphi \leq 2\pi. \quad (5)$$

### B. Fixing spin

sec:theory-fixspin

In a similar way we can fix in the  $xy$  plane the spin direction along the  $x$ ,  $y$ , and  $z$  directions and then define the magnitude of the spin velocity  $|\mathcal{V}_{\sigma^b}|$  in a fixed angle  $\gamma_b$

$$|\mathcal{V}_{\sigma^b}| = \sqrt{(\mathcal{V}^{ax})^2 + (\mathcal{V}^{ay})^2}, \quad (6)$$

$$\gamma_b = \tan^{-1} \left( \frac{\mathcal{V}^{ay}}{\mathcal{V}^{ax}} \right), \quad (7)$$

where the angle is measured in the counter-clockwise direction from the positive  $x$  axis.

## III. RESULTS

sec:results

We preset the results for  $\mathcal{V}^{ab}$  for the  $\text{C}_{16}\text{H}_8$ -alt and  $\text{C}_{16}\text{H}_8$ -up structures being both noncentrosymmetric semi-infinite carbon systems with 50% hydrogenation in different arrangements. The *alt* system has alternating hydrogen atoms on the upper and bottom sides of the carbon sheet, while the *up* system has H only on the upper side. We take the hexagonal carbon lattice to be on the  $xy$  plane for both structures, and the carbon-hydrogen bonds on the perpendicular  $xz$  plane, as depicted in Figs. 1 and 2.

Using the ABINIT code<sup>1</sup> we calculated the self-consistent ground state and the Kohn-Sham states using density functional theory in the local density approximation (DFT-LDA) with a planewave basis. We used Hartwigsen-

Layer No.	Atom type	Position [ $\text{\AA}$ ]		
		$x$	$y$	$z$
1	H	-0.61516	-1.42140	1.47237
2	C	-0.61516	-1.73300	0.39631
3	C	0.61516	1.73300	0.15807
4	C	0.61516	0.42201	-0.15814
5	C	-0.61516	-0.37396	-0.39632
6	H	-0.61516	-0.68566	-1.47237

TABLE I. Unit cell of *alt* structure. Layer division, atom types and positions for the *alt* structure. The structure unit cell was divided in six layers corresponding each one to atoms in different  $z$  positions. The corresponding layer atom position is depicted in Fig. 1 with the corresponding number of layer.

tab:alt-unitcell

Goedecker-Hutter (HGH) relativistic separable dual-space Gaussian pseudopotentials<sup>2</sup> including the spin-orbit interaction for calculating  $\mathcal{V}^a(\omega)$ .

The convergence parameters for the calculations of our results corresponding to the *alt* and *up* structures are cutoff energies of 65 Ha and 40 Ha, respectively. The energy eigenvalues and matrix elements were calculated using 14452  $\mathbf{k}$  points and 8452  $\mathbf{k}$  points in the irreducible Brillouin zone (IBZ) and present LDA energy band gaps of 0.72 eV and 0.088 eV, respectively for the *alt* and *up* structures. As mentioned in<sup>3</sup>, using DFT the LDA is only one method of many other that can be used to calculate the electronic structure of materials. Also it is known that all methods predict a different band gap than the obtained in the experiment. A correction for the band gap energy value can be calculated by other *ab-initio* methods such as the GW approximation<sup>4</sup> being this outside the scope of this paper.

The structures presented here were divided into layers to analyze the layer-by-layer contribution for  $\mathcal{V}^{ab}$  response. The *alt* structure was divided in six layers corresponding the first one to the top hydrogen atoms, from the second to the forth to carbon atoms in different  $z$  positions, and the sixth and last one to the bottom hydrogen atoms. The *up* structure was divided into two layers, the first one comprised by the top hydrogen atoms and the second by the carbon atoms. The layer divisions and atom positions

Layer No.	Atom type	Position [ $\text{\AA}$ ]		
		$x$	$y$	$z$
1	H	-0.61516	-1.77416	0.73196
1	H	0.61518	0.35514	0.73175
2	C	-0.61516	-1.77264	-0.49138
2	C	-0.61516	-0.35600	-0.72316
2	C	0.61516	0.35763	-0.49087

TABLE II. Unit cell of *up* structure. Layer division, atom types and positions for the *up* structure. The structure unit cell was divided in two layers corresponding to hydrogen and carbon atoms. The corresponding layer atom position is depicted in Fig. 2 with the corresponding number of layer.

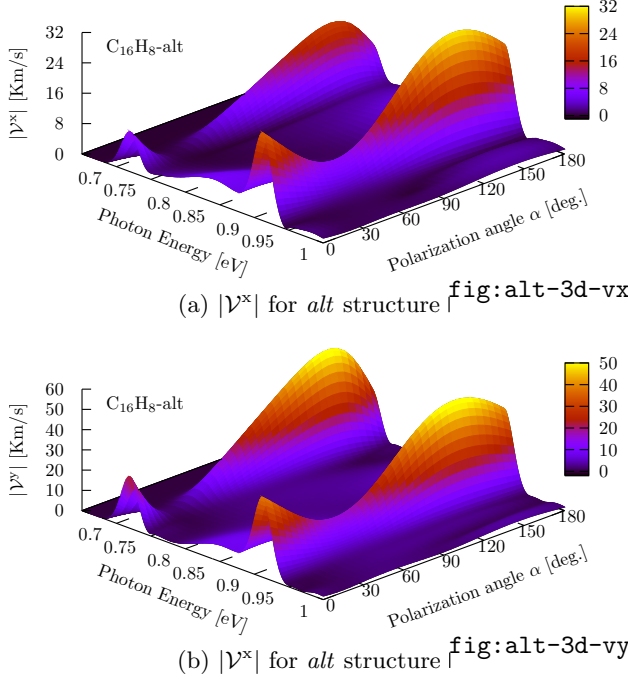


FIG. 3.  $|V^x|$  response for  $C_{16}H_8$ -*alt* structure. The maximum response zone is localized for an energy range from 0.90 eV to 0.93 eV,  $145^\circ$  and for a polarization angle of the incoming beam from  $120^\circ$  to  $150^\circ$ .

for the unit cells are shown in Tables I and II.

### A. Fixing velocity

sec:res-fixvel

For the *alt* structure we analyzed the energy range of energy from 0.6 eV to 1.0 eV where we found the most intense response for  $V^{ab}$  and  $|V^a|$ .

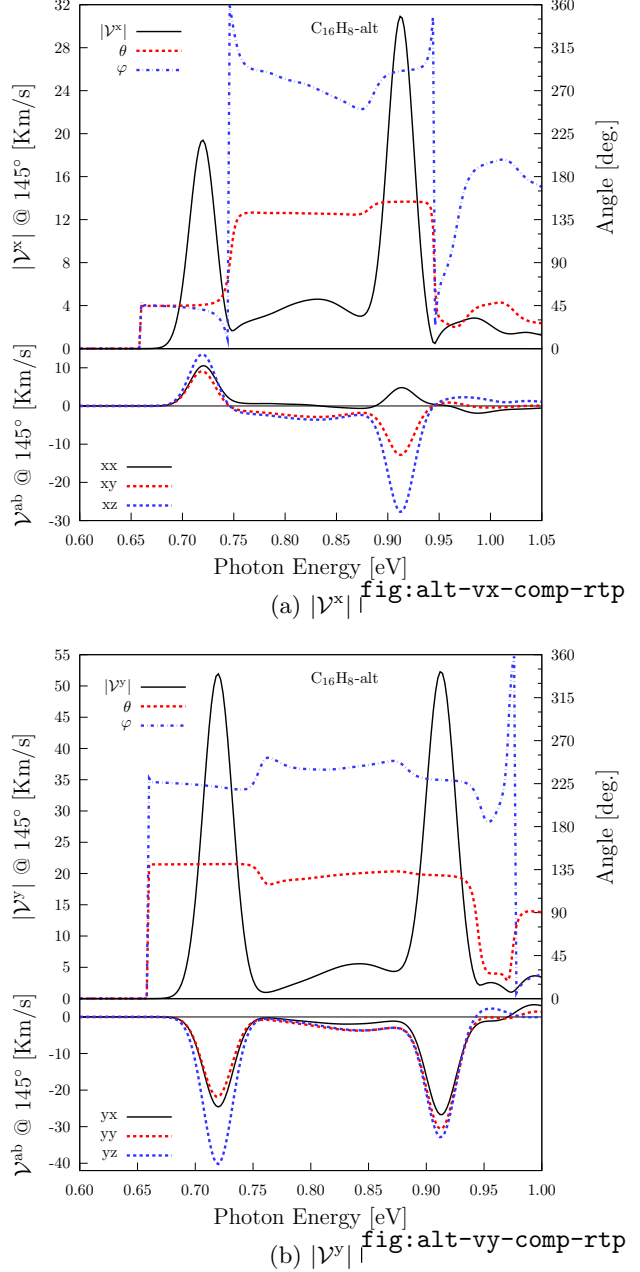


FIG. 4. Most intense responses of  $|V^x|$  and  $|V^y|$  and the corresponding three components for the *alt* structure. Both maxima were obtained for a polarization angle  $\alpha = 145^\circ$ .

In Fig. 3 we present the  $|V^a|$  spectra resulting from evaluate Eq. (3) using different polarization angles  $\alpha$  in Eq. (1) for the  $C_{16}H_8$ -*alt* structure. We can see that the onset of the response is when the energy of the incoming light is the same of the gap energy. From this picture we can see that for the zone between the

energy range of 0.90 eV-0.93 eV and polarization angles between  $120^\circ$  and  $150^\circ$  is the zone of the absolute maximum response for both,  $|\mathcal{V}^x|$  and  $|\mathcal{V}^y|$ . Also there is another zone of interest for energies from 0.70 eV to 0.74 eV where a local maximum is obtained. From Fig. 3(a) we have that  $|\mathcal{V}^x|$  reaches values of 30 Km/s for the first zone mentioned before and 20 Km/s for the second one. We also found that the absolute maximum of the response is obtained when the polarization angle is  $\alpha = 145^\circ$ . In the top frames of Figs. 4(a) and 4(b) we present the results for  $|\mathcal{V}^x|$  and  $|\mathcal{V}^y|$  fixing the polarization angle to  $145^\circ$  for the *alt* structure vs the photon energy and the corresponding azimuthal  $\theta$  and polar  $\varphi$  angles. Also in the bottom frames of Figs. 4(a) and 4(b) we present the decomposition of  $|\mathcal{V}^x|$  and  $|\mathcal{V}^y|$  in the corresponding  $\mathcal{V}^{xx}$ ,  $\mathcal{V}^{xy}$ ,  $\mathcal{V}^{xz}$  and  $\mathcal{V}^{yx}$ ,  $\mathcal{V}^{yy}$ ,  $\mathcal{V}^{yz}$  components for the fixed polarization angle. Making the analysis for the components and angles for  $|\mathcal{V}^x|$  depicted in Fig. 4(a) we can see that for the energy range from 0.70 eV to 0.74 eV all the  $xx$ ,  $xy$ , and  $xz$  components contribute with almost the same intensity giving a total spin-velocity of 19.3 Km/s and spin polar and azimuthal angles  $\varphi = 45.8^\circ$  and  $\theta = 40.7^\circ$ . In the other hand, for the energy range from 0.88 eV to 0.95 eV there is a major contribution coming from the  $\mathcal{V}^{xz}$  component resulting in a spin-velocity magnitude of 30.9 Km/s being this magnitude the most intense for  $|\mathcal{V}^x|$ . In this case the polar angle is  $\varphi = 153.8^\circ$  and the spin angle over the  $xy$  plane have is  $\theta = 290.4^\circ$ . Also we notice that for the range of 0.70-0.74 eV all the contributions are positive while for the range of 0.88-0.95 eV the  $xx$  component remains positive but the components  $xy$  and  $xz$  change in direction. This is due to **a change in the spin polarization**. Making now the analysis for  $|\mathcal{V}^y|$  depicted in Fig. 4(b) we have that for the energy range from 0.70 eV to 0.74 eV the  $yz$  component have a more intense response than  $yx$  and  $yy$  components and they give a total spin-velocity of 51.9 Km/s and result in polar angle  $\varphi = 140.7^\circ$  and spin azimuthal angle of  $\theta = 221.5^\circ$ . For the energy range from 0.88 eV to 0.95 eV all three components have similar intensities resulting in a spin-velocity magnitude of  $|\mathcal{V}^y| = 52.3$  Km/s being this response 1.7 times mores intense than

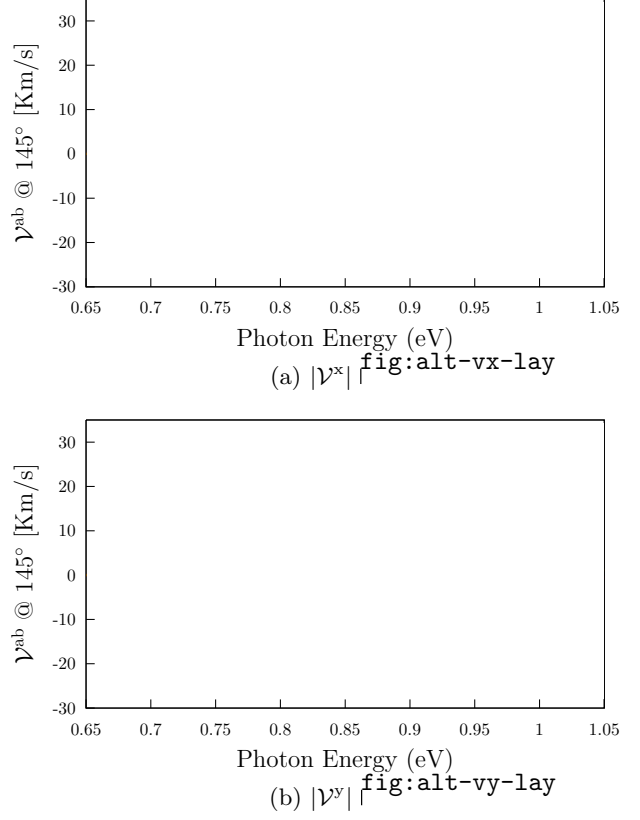


FIG. 5. Layer-by-layer contribution of  $|\mathcal{V}^x|$  and  $|\mathcal{V}^y|$  for a polarization angle  $\alpha = 145^\circ$  for the *alt* structure.

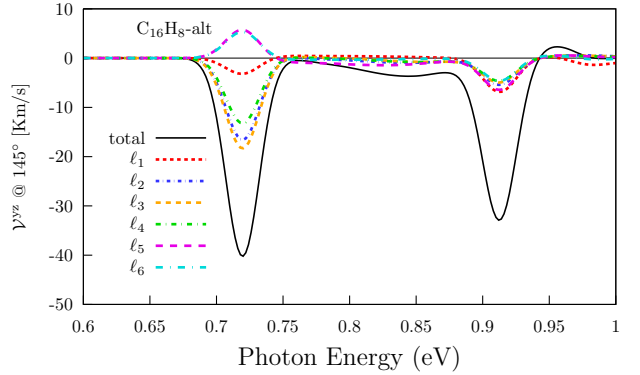


FIG. 6. Layer-by-layer contribution of  $\mathcal{V}^{yz}$  for the *alt* structure.

the most intense response of  $|\mathcal{V}^x|$ . The corresponding polar and azimuthal angles for  $|\mathcal{V}^y|$  in this energy range are  $\theta = 129.0^\circ$  and  $\varphi = 228.9^\circ$ . Now we have that the three components of  $|\mathcal{V}^y|$  are negative for the energy range from 0.60 eV to 1.0 eV **keeping the same spin polarization for all this range**.

In Fig. 5 we show the layer-by-layer contri-

bution of  $|\mathcal{V}^{ab}|$  for the *alt* structure. The corresponding layer division and atom types and positions are presented in Table I and depicted in Fig. 1. In the layer- by-layer contribution for  $|\mathcal{V}^x|$  (Fig. 5(a)) we have that ... Also, for the layer-by-layer contribution for  $|\mathcal{V}^y|$  (Fig. 5(b)) we have that ... From the bottom panels of Fig. 4 we can see that for the *alt* structure the most intense component of  $|\mathcal{V}^x|$  and  $|\mathcal{V}^y|$  corresponds to  $\mathcal{V}^{yz}$  which has a value of -40.21374498 Km/s for an energy incident beam of 0.72 eV. This component and the corresponding layer by layer contribution is depicted in Fig. 6. From this figure we have that for the energy range from 0.70 eV to 0.74 eV the fifth and sixth layers cor-

responding to the bottom carbon and hydrogen numbered with 5 and 6 in Fig. 1 have contributions in opposite direction than the other 4 layers resulting in a total response  $\mathcal{V}^{yz} = -40.2$  Km/s for an incoming beam energy of 0.72 eV. In the other hand, for the energy range from 0.88 eV to 0.95 eV the response for the all six layers the responses are in the same direction resulting in a total response  $\mathcal{V}^{yz} = -32.89$  Km/s for an incoming beam with energy of 0.912 eV.

## B. Fixing spin

sec:res-fixspin

---

<sup>1</sup> X. Gonze, B. Amadon, P.-M. Anglade, J.-M. Beuken, F. Bottin, P. Boulanger, F. Bruneval, D. Caliste, R. Caracas, M. Côté, T. Deutsch, L. Genovese, P. Ghosez, M. Giantomassi, S. Goedecker, D. Hamann, P. Hermet, F. Jollet, G. Jomard, S. Leroux, M. Mancini, S. Mazevet, M. Oliveira, G. Onida, Y. Pouillon, T. Rangel, G.-M. Rignanese, D. Sangalli, R. Shaltaf, M. Torrent, M. Verstraete, G. Zerah, and J. Zwanziger, Com-

put. Phys. Commun. **180**, 2582 (2009).

<sup>2</sup> C. Hartwigsen, S. Goedecker, and J. Hutter, Phys. Rev. B **58**, 3641 (1998).

<sup>3</sup> R. Zapata-Peña, S. M. Anderson, B. S. Mendoza, and A. I. Shkrebtii, physica status solidi (b) **253**, 226 (2016).

<sup>4</sup> G. Onida, L. Reining, and A. Rubio, Rev. Mod. Phys. **74**, 601 (2002)

LastBibItem