



Nombres:

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Título Profesional o Grado Académico (incluya el año de obtención):

**LICENCIADO EN FÍSICA, UNIVERSIDAD DE LA HABANA ,1995.**

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Estudios de Postgrado o Especialización (institución donde lo obtuvo y año de obtención):

**MÁSTER EN CIENCIAS FÍSICAS, UNIVERSIDAD DE LA HABANA, CUBA, 1997.**

**DOCTOR EN CIENCIAS FÍSICAS, UNIVERSIDAD DE LA HABANA, CUBA, 2001**

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Actividad Actual e Institución en la cual trabaja:

**PROFESOR ASISTENTE DE LA UNIVERSIDAD DE CHILE. DEPARTAMENTO DE FÍSICA, FACULTAD DE CIENCIAS.**

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Reseña de su actividad laboral actual:

**Docencia:** Imparto Física en las carreras de la Facultad. También doy cursos de postgrado del programa de doctorado en física. Asimismo, trabajo en proyectos de investigación en ciencia de materiales.

**Líneas de Investigación:** estoy interesado en las aplicaciones de la simulación computacional de la escala atómica a la predicción y explicación de las propiedades de los materiales. Principalmente aquellos problemas donde la combinación de la experimentación y la simulación es necesarias para obtener respuestas. Algunos ejemplos son la espectroscopia de los materiales complejos y nanoestructuras, la estabilidad de fase en condiciones extremas, los materiales amorfos y poco cristalino,

y materiales promisorios con métodos de fabricación aún no desarrollados. Tengo experiencia en los cálculos de estructura electrónica, dinámica molecular, propiedades elásticas, fonones, espectros ópticos y espectros de niveles profundos. La mayoría de mis cálculos de estructura electrónica se basan en la teoría funcional de la densidad, pero también utilizan métodos de función de onda. En particular, actualmente estoy trabajando en la aplicación y desarrollo de un método de función de onda aproximada para la descripción de las excitaciones electrónicas en los sistemas moleculares con gran número de átomos. También uso, siempre que sea posible, las técnicas de simulación basadas en potenciales interatómicos empíricos.

Información adicional: <http://fisica.ciencias.uchile.cl/~emenendez>

#### PUBLICACIONES:

Montero-Alejo, A.L., Menéndez-Proupin, E., Fuentes, M.E., Delgado, A., Montforts, F.-P., Montero-Cabrera, L.A., García De La Vega, J.M.  
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Menéndez-Proupin, E., Montero-Alejo, A.L., García De La Vega, J.M.  
Ultrathin carbon nanotube with single, double, and triple bonds  
(2012) Physical Review Letters, 109 (10), art. no. 105501, .  
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Margapoti, E., Alves, F.M., Mahapatra, S., Lopez-Richard, V., Worschech, L., Brunner, K., Qu, F., Destefani, C., Menéndez-Proupin, E., Bougerol, C., Forchel, A., Marques, G.E.  
Paramagnetic shift in thermally annealed Cd xZn 1-xSe quantum dots  
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Menéndez-Proupin, E., Cervantes-Rodríguez, S., Osorio-Pulgar, R., Franco-Cisterna, M., Camacho-Montes, H., Fuentes, M.E.  
Computer simulation of elastic constants of hydroxyapatite and fluorapatite  
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Ortega, Y., Hernández, N.C., Menéndez-Proupin, E., Graciani, J., Sanz, J.F.  
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SOURCE: Scopus

Margapoti, E., Alves, F.M., Mahapatra, S., Schmidt, T., Lopez-Richard, V., Destefani, C., Menéndez-Proupin, E., Qu, F., Bougerol, C., Brunner, K., Forchel, A., Marques, G.E., Worschech, L.  
Characterization of spin-state tuning in thermally annealed semiconductor quantum dots  
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Lizárraga, R., Holmström, E., Amézaga, A., Bock, N., Peery, T., Menéndez-Proupin, E., Giannozzi, P.  
Core-level shift analysis of amorphous CdTeO<sub>x</sub> materials  
(2010) Journal of Materials Science, 45 (18), pp. 5071-5076. Cited 4 times.  
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Loyola, C., Menéndez-Proupin, E., Gutiérrez, G.  
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Montero-Alejo, A.L., Fuentes, M.E., Menéndez-Proupin, E., Orellana, W., Bunge, C.F., Montero, L.A., García De La Vega, J.M.

Approximate quantum mechanical method for describing excitations and related properties of finite single-walled carbon nanotubes  
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Ab initio molecular dynamics study of amorphous CdTeOx alloys: Structural properties  
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Exciton-phonon complexes and optical properties in CdSe nanocrystals  
(2006) Journal of Physics Condensed Matter, 18 (31), art. no. 022, pp. 7283-7298.  
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Menéndez-Proupin, E.  
Comment on "The effects of electric field on the electronic structure of a semiconductor quantum dot" [J. Appl. Phys. 84, 1454 (1998)]  
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<http://www.scopus.com/inward/record.url?eid=2-s2.0-2342440425&partnerID=40&md5=d4a8242b0cc6380292870569583a2db8>  
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Menéndez-Proupin, E., Cabo-Bisset, N.  
Resonance Raman scattering in semiconductor quantum dots: Adiabatic versus time-dependent perturbation theory  
(2002) *Physical Review B - Condensed Matter and Materials Physics*, 66 (8), art. no. 085317, pp. 853171-853178. Cited 4 times.  
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Resonant Raman scattering off neutral quantum dots  
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Iribarren, A., Menéndez-Proupin, E., Caballero-Briones, F., Castro-Rodríguez, R., Peña, J.L.  
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(2001) *Modern Physics Letters B*, 15 (17-19), pp. 643-646. Cited 4 times.  
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Gonzalez, A., Menéndez-Proupin, E.  
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Rodríguez-Suárez, R., Menéndez-Proupin, E., Trallero-Giner, C., Cardona, M.  
Multiphonon resonant Raman scattering in nanocrystals  
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Menéndez-Proupin, E., Trallero-Giner, C., Ulloa, S.E.  
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**PROYECTOS DE INVESTIGACIÓN:**

Investigador Responsable. 1120183 Study of impurities, complex defects, and grain boundaries in cdte thin films using computer simulation at the atomic scale. 2012

Investigador Responsable. 7060248 Physics of novel solar cell materials. 2006

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Investigador extranjero de contraparte. 7030083 Study of properties of materials by computer simulation: from atomic to nanoscopic regime. Estudio de la propiedades de materiales mediante simulación computacional: de lo atómico a lo nanoscópico. 2003

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