



Nombres:

FERNANDO JAVIER

Apellidos:

MENDIZÁBAL EMALDÍA

Contacto (Opcional):

HAGUA@UCHILE.CL

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

LICENCIATURA EN CIENCIAS C/M EN QUÍMICA, UNIVERSIDAD DE CHILE 1990.

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

**DOCTOR EN CIENCIAS C/M EN QUÍMICA, UNIVERSIDAD DE CHILE 1995;
ESTADÍA POSTDOCTORAL U. DE HELSINKI (FINLANDIA) EN QUÍMICA
INORGÁNICA TEÓRICA, 1996-1997.**

Actividad Actual e Institución en la cual trabaja:

**PROFESOR TITULAR DE LA UNIVERSIDAD DE CHILE. DIRECTOR DEL
DEPARTAMENTO DE QUÍMICA DE LA FACULTAD DE CIENCIAS Y
CONSEJERO DE FACULTAD.**

Reseña de su actividad laboral actual:

Área de Investigación: Química inorgánica teórica y computacional.

Línea de Investigación: Físicoquímica inorgánica teórica de interacciones débiles de Van der Waals en moléculas y sólidos extendidos de metales pesados de oro, mercurio, talio, etc; polímeros inorgánicos de coordinación basados en metalomacrociclos puenteados con ligandos nitrogenados biaxiales; índices de

reactividad química en los esquemas ab initio y funcionales de la densidad en química de materiales y clusters metálicos.

PUBLICACIONES INDEXADAS:

Mendizabal, F., Salazar, R.

Theoretical study on electronic spectra and interaction in $[Au_3]-L-[Au_3]$ ($L = C_6F_6, Ag^+$) complexes

(2012) Journal of Molecular Modeling, pp. 1-7. Article in Press.

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DOCUMENT TYPE: Article in Press

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Ponce, I., Silva, J.F., Oñate, R., Rezende, M.C., Paez, M.A., Zagal, J.H., Pavez, J., Mendizabal, F., Miranda-Rojas, S., Muñoz-Castro, A., Arratia-Pérez, R.

Enhancement of the catalytic activity of Fe phthalocyanine for the reduction of O_2 anchored to Au(111) via conjugated self-assembled monolayers of aromatic thiols as compared to Cu phthalocyanine

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Aguilera-Venegas, B., Olea-Azar, C., Arán, V.J., Maya, J.D., Kemmerling, U., Speisky, H., Mendizabal, F.

Electrochemical, ESR and theoretical insights into the free radical generation by 1,1'-hydrocarbylenebisindazoles and its evaluation as potential bio-active compounds

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Mera-Adasme, R., Mendizabal, F., Gonzalez, M., Miranda-Rojas, S., Olea-Azar, C., Sundholm, D.

Computational studies of the metal-binding site of the wild-type and the H46R mutant of the copper, zinc superoxide dismutase

(2012) Inorganic Chemistry, 51 (10), pp. 5561-5568.

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Burgos, D., Olea-Azar, C., Mendizabal, F.

Theoretical study of the local reactivity of electrophiles of the type M^+PR_3 ($M=Cu, Ag, Au$; $R=H, -Me, -Ph$)

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Theoretical and experimental study of bonding and optical properties of self-assembly metallophthalocyanines complexes on a gold surface. A survey of the substrate-surface interaction

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Basicity of bisperhalophenyl aurates toward closed-shell metal ions: Metallophilicity and additional interactions

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Theoretical study of the interaction between Pt(0) and MPH₃ + fragments in complexes of the $[\text{Pt}3(\mu\text{-CO})_3(\text{PH}_3)_3\text{-MPH}_3 + (\text{M} = \text{Cu}^+, \text{Au}^+, \text{Ag}^+)]$ type

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Mera-Adasme, R., Mendizabal, F., Olea-Azar, C., Miranda-Rojas, S., Fuentealba, P.
A computationally efficient and reliable bond order measure

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Rodríguez-Castillo, M., Monge, M., López-de-Luzuriaga, J.M., Olmos, M.E., Laguna, A., Mendizabal, F.

Theoretical study of the closed-shell d₁₀-d₁₀ Au(I)-Cu(I) attraction in complexes in extended unsupported chains

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Theoretical study of the s₂-p interaction between t_l(i) and olefinic systems on hypothetical [t_l(c₂h₄)₂]⁺ and [t_l(c₂h₂)_n]⁺ (n= 2,3) complexes
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Theoretical EPR study of nitroindazoles: 3-alkoxy, 3-hydroxy and 3-oxo derivatives
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Theoretical study of {au₃(ch₃nacoch₃)₃}_n.{2,4,7-trinitro-9-?uorenone} (n = 1,2) complexes
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Mera, R., Mendizabal, F.

Theoretical study of the d10-s2 interaction between Au(I) and Tl(I) in the [AuCl(PH₃)₂]Tl⁺ hypothetical complex
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Mendizabal, F., Burgos, D., Olea-Azar, C.
Theoretical study of electronic spectra of [Pt₃(μ-CO) ₃(CO)₃]_n -2 (n = 3-5) complexes
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Jullian, C., Orosteguis, T., Pérez-Cruz, F., Sánchez, P., Mendizabal, F., Olea-Azar, C.
Complexation of morin with three kinds of cyclodextrin. A thermodynamic and reactivity study
(2008) Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 71 (1), pp. 269-275. Cited 34 times.
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Mendizabal, F., Burgos, D., Olea-Azar, C.
Theoretical study of [Hg₃(o-C₆F₄)₃]_n · {benzene} (n = 1, 2) complexes
(2008) Chemical Physics Letters, 463 (1-3), pp. 272-277. Cited 4 times.
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Fernández, E.J., Laguna, A., López-de-Luzuriaga, J.M., Monge, M., Mendizabal, F.
Theoretical study of the aggregation of d10-s2 Au(I)-Tl(I) complexes in extended unsupported chains
(2008) Journal of Molecular Structure: THEOCHEM, 851 (1-3), pp. 121-126. Cited 5 times.
<http://www.scopus.com/inward/record.url?eid=2-s2.0-39149119965&partnerID=40&md5=c29a8002e0a02d83dc584552e9b7b582>
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Theoretical study of the electronic spectra of Bi- and Tri-heteronuclear platinum complexes
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Theoretical study on electronic spectra and aurophilic attraction in $[\text{Au}_3(\text{MeN}\{\text{double bond, long}\}\text{COMe})_3]_n$ ($n = 1$ s(-) 4) complexes

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Jullian, C., Miranda, S., Zapata-Torres, G., Mendizábal, F., Olea-Azar, C.

Studies of inclusion complexes of natural and modified cyclodextrin with (+)catechin by NMR and molecular modeling

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Mendizábal, F., Olea-Azar, C., Miranda, S.

Theoretical study of the interaction d10-d8 between Pt(0) and M(I) on the $[\text{Pt}(\text{PH}_3)_3\text{MPH}_3]^+$ complexes ($M = \text{Cu}, \text{Ag}, \text{Au}$)

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Mendizabal, F., Donoso, D., Olea-Azar, C., Mera, R.

Theoretical study of the interaction d10-s2 between Pt(0) and metals on the $[\text{Pt}(\text{PH}_3)_3\text{M}]$ complexes ($M = \text{Hg}(0), \text{Au}(-\text{I})$)

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Mendizabal, F., Olea-Azar, C.

Theoretical study in $[\text{C}_2\text{H}_4\text{-Ti}]^+$ and $[\text{C}_2\text{H}_2\text{-Ti}]^+$ complexes

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Rigol, C., Olea-Azar, C., Mendizábal, F., Briones, R., Cerecetto, H., González, M.

Theoretical study on 5-nitrofuryl thiosemicarbazone radicals electronic properties

(2006) Journal of Molecular Structure: THEOCHEM, 770 (1-3), pp. 125-129. Cited 4 times.

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Theoretical study on the electronic spectrum of $[M(CN)_2]_n - n$ ($M=Au(I), Ag(I); n=1-3$) complexes
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Electronic structure and molecular properties of the octacyanorhenate $[Re(CN)_8]^{3-}$ and $[Re(CN)_8]^{2-}$ complexes
(2006) Chemical Physics Letters, 422 (1-3), pp. 89-94. Cited 2 times.
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Complexes self-associate by hydrogen bonding and metallophilic attraction: Theoretical study
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Olea-Azar, C., Rigol, C., Mendizábal, F., Briones, R.
Applications of electron spin resonance and spin trapping in tropical parasitic diseases
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Rigol, C., Olea-Azar, C., Mendizábal, F., Otero, L., Gambino, D., González, M., Cerecetto, H.
Electrochemical and ESR study of 5-nitrofuryl-containing thiosemicarbazones antiprotozoal drugs
(2005) Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 61 (13-14), pp. 2933-2938. Cited 23 times.
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