

# Eberth Correa

## List of Publications by Year in descending order

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Version: 2024-02-01

22  
papers

93  
citations

1937685  
4  
h-index

1474206  
9  
g-index

23  
all docs

23  
docs citations

23  
times ranked

51  
citing authors

#	ARTICLE	IF	CITATIONS
1	Field-theoretical renormalization group for a flat two-dimensional Fermi surface. Physical Review B, 2005, 71, .	3.2	32
2	Breakdown of the Fermi-liquid regime in the two-dimensional Hubbard model from a two-loop field-theoretical renormalization group approach. Physical Review B, 2008, 78, .	3.2	16
3	Quantum and classical study of vibrational states of H <sub>2</sub> and H <sub>3</sub> molecules. International Journal of Quantum Chemistry, 2003, 95, 149-152.	2.0	6
4	Two-loop renormalization group calculation of response functions for a two-dimensional flat Fermi surface. Physical Review B, 2008, 78, .	3.2	5
5	Spherical harmonics representation of the potential energy surface for the H <sub>2</sub> -H <sub>2</sub> van der Waals complex. Journal of Molecular Modeling, 2020, 26, 277.	1.8	5
6	Theoretical study of the H + HCN → H + HNC process. Journal of Molecular Modeling, 2017, 23, 169.	1.8	4
7	Renormalization group calculation of the uniform susceptibilities in low-dimensional systems. Journal of Physics A, 2006, 39, 7977-7992.	1.6	3
8	Insulating spin liquid in the lightly doped two-dimensional Hubbard model. Physical Review B, 2006, 73, .	3.2	3
9	A Functional Generalization of the Field-Theoretical Renormalization Group Approach for the Single-Impurity Anderson Model. Journal of Low Temperature Physics, 2012, 166, 192-207.	1.4	3
10	Fermi surface renormalization in two dimensions. Physica C: Superconductivity and Its Applications, 2004, 408-410, 254-256.	1.2	2
11	SUPERCONDUCTIVITY IN THE 2D ATTRACTIVE HUBBARD MODEL WITHIN A FUNCTIONAL FIELD-THEORETICAL RG. International Journal of Modern Physics B, 2011, 25, 3691-3706.	2.0	2
12	MOLECULAR DYNAMICS SIMULATION OF A TWO-DIMENSIONAL HEISENBERG FLUID. International Journal of Modern Physics C, 2012, 23, 1250026.	1.7	2
13	Rate constant calculations of the C <sub>2</sub> + HCN → CCCN+H addition via the Master Equation. Journal of Molecular Modeling, 2017, 23, 143.	1.8	2
14	Thermal rate constant for the C <sub>3</sub> P + OH(X <sub>2</sub> I) → CO(X <sub>1</sub> I) + H <sub>2</sub> S reaction using stochastic energy grained master equation method. International Journal of Chemical Kinetics, 2019, 51, 590-601.	1.6	2
15	Isocyanoacetylene and cyanoacetylene formation study from C <sub>2</sub> H <sub>2</sub> + CN reaction. International Journal of Chemical Kinetics, 2022, 54, 309-316.	1.6	2
16	Oscilador harmônico com massa variável e a segunda lei de Newton. Revista Brasileira De Ensino De Fisica, 2011, 33, 4307-4307.	0.2	1
17	Fermi surface renormalization and quantum confinement in the two-coupled chains model. European Physical Journal B, 2014, 87, 1.	1.5	1
18	Kinetics and mechanism of the $\text{CH}_3\text{OH} + \text{H}_2\text{S} \rightarrow \text{CH}_3\text{SH} + \text{H}_2\text{O}$ reaction. Chemical Physics Letters, 2019, 734, 136699.	2.6	1

#	ARTICLE	IF	CITATIONS
19	Long range strength and anisotropies of molecule-molecule interactions: Ab initio calculations, spherical harmonics expansions, and the second virial coefficient for the $\text{H}_2\text{H}_2$ system	2.6	1
20	Graphical Visualization on Computational Simulation Using Shared Memory. Journal of Physics: Conference Series, 2014, 487, 012014.	0.4	0
21	Solution of the 1d Schrödinger Equation for a Symmetric Well. Revista Brasileira De Ensino De Fisica, 2019, 41, .	0.2	0
22	A Simple Monte Carlo Simulation For the Two Dimensional Attractive Hubbard Model. Journal of Physics: Conference Series, 2020, 1483, 012002.	0.4	0