

Juan M Sánchez

List of Publications by Year in descending order

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104
docs citations

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times ranked

2766
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | Generalized cluster description of multicomponent systems. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1984, 128, 334-350. | 2.6 | 1,536 |
| 2 | The fee Ising model in the cluster variation approximation. <i>Physical Review B</i> , 1978, 17, 2926-2936. | 3.2 | 336 |
| 3 | Ordering in fcc lattices with first- and second-neighbor interactions. <i>Physical Review B</i> , 1980, 21, 216-228. | 3.2 | 200 |
| 4 | Local stability of nonequilibrium phases. <i>Physical Review Letters</i> , 1994, 72, 3076-3079. | 7.8 | 164 |
| 5 | Ni-Pt Phase Diagram: Experiment and Theory. <i>Physical Review Letters</i> , 1985, 55, 1208-1211. | 7.8 | 161 |
| 6 | Overview no. 43. <i>Acta Metallurgica</i> , 1985, 33, 1171-1185. | 2.1 | 142 |
| 7 | First-principles calculation of the Ag-Cu phase diagram. <i>Physical Review B</i> , 1991, 44, 5411-5418. | 3.2 | 135 |
| 8 | Theoretical calculation of the Cu ₃ -Ag-Au coherent phase diagram. <i>Acta Metallurgica</i> , 1980, 28, 651-662. | 2.1 | 127 |
| 9 | Cluster expansion and the configurational theory of alloys. <i>Physical Review B</i> , 2010, 81, . | 3.2 | 126 |
| 10 | Comparison of approximate methods for the study of antiferromagnetism in the fcc lattice. <i>Physical Review B</i> , 1982, 26, 1465-1468. | 3.2 | 124 |
| 11 | Density-functional study of Fe ₃ Al: LSDA versus GGA. <i>Physical Review B</i> , 2002, 65, . | 3.2 | 119 |
| 12 | Modeling of $\hat{\Gamma}_3/\hat{\Gamma}_2$ phase equilibrium in the nickel-aluminum system. <i>Acta Metallurgica</i> , 1984, 32, 1519-1525. | 2.1 | 113 |
| 13 | Cluster expansions and the configurational energy of alloys. <i>Physical Review B</i> , 1993, 48, 14013-14015. | 3.2 | 113 |
| 14 | Calculation of phase equilibrium in Al-Li alloys. <i>Acta Metallurgica</i> , 1986, 34, 1021-1028. | 2.1 | 106 |
| 15 | Ising model phase-diagram calculations in the fcc lattice with first- and second-neighbor interactions. <i>Physical Review B</i> , 1982, 25, 1759-1765. | 3.2 | 97 |
| 16 | Model for Anomalous Self-Diffusion in Group-IVB Transition Metals. <i>Physical Review Letters</i> , 1975, 35, 227-230. | 7.8 | 95 |
| 17 | Short range order diffuse intensity calculations in the cluster variation method. <i>Acta Metallurgica</i> , 1985, 33, 1463-1474. | 2.1 | 92 |
| 18 | Magnetic properties and chemical ordering in Co-Pt. <i>Journal of Physics Condensed Matter</i> , 1989, 1, 491-496. | 1.8 | 88 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 19 | Theoretical description of phase equilibrium in binary alloys. <i>Acta Metallurgica</i> , 1985, 33, 1097-1104. | 2.1 | 87 |
| 20 | Structure and magnetism in bcc-based iron-cobalt alloys. <i>Physical Review B</i> , 2006, 73, . | 3.2 | 87 |
| 21 | First-principles study of the solubility of Zr in Al. <i>Physical Review B</i> , 2002, 65, . | 3.2 | 84 |
| 22 | Dynamical Scaling in the Glass System $B_2O_3-PbO-Al_2O_3$. <i>Physical Review Letters</i> , 1981, 47, 1308-1311. | 7.8 | 82 |
| 23 | Structural instabilities of excited phases. <i>Physical Review B</i> , 1997, 55, 787-797. | 3.2 | 74 |
| 24 | First-principles investigation of the Ni-Fe-Al system. <i>Intermetallics</i> , 2005, 13, 1096-1109. | 3.9 | 73 |
| 25 | Cluster-method study of surface effects in fcc binary alloys. <i>Physical Review B</i> , 1985, 32, 3534-3540. | 3.2 | 71 |
| 26 | Ordering and segregation at (001) surfaces of Cu_3Au . <i>Surface Science</i> , 1985, 157, L297-L302. | 1.9 | 70 |
| 27 | Anomalous diffusion in omega forming systems. <i>Acta Metallurgica</i> , 1978, 26, 1083-1095. | 2.1 | 64 |
| 28 | First-order phase transitions in the Ising square lattice with first- and second-neighbor interactions. <i>Physical Review B</i> , 1993, 48, 3519-3522. | 3.2 | 62 |
| 29 | Modeling of ternary site occupation in L12 ordered intermetallics. <i>Acta Metallurgica</i> , 1989, 37, 2835-2840. | 2.1 | 61 |
| 30 | Pair correlations in the cluster variation approximation. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1982, 111, 200-216. | 2.6 | 60 |
| 31 | New evidence for GP zones in binary Al-Li alloys. <i>Scripta Metallurgica</i> , 1986, 20, 201-206. | 1.2 | 60 |
| 32 | Calculation of Thermodynamic Properties and Phase Diagrams of Binary Transition-Metal Alloys. <i>Physical Review Letters</i> , 1986, 57, 253-256. | 7.8 | 58 |
| 33 | Phase diagram and thermodynamic properties of Ni-Al alloys: A non-empirical evaluation. <i>Solid State Communications</i> , 1988, 65, 527-530. | 1.9 | 52 |
| 34 | First-principles study of short range order and instabilities in AuCu, AuAg and AuPd alloys. <i>Acta Metallurgica Et Materialia</i> , 1991, 39, 493-501. | 1.8 | 51 |
| 35 | Spontaneous L12 Order at Ni90Al10(110) Surfaces: An X-Ray and First-Principles-Calculation Study. <i>Physical Review Letters</i> , 2001, 87, 236102. | 7.8 | 51 |
| 36 | Ordering and layer composition at the $Cu_3Au(110)$ surface. <i>Surface Science</i> , 1990, 238, L481-L485. | 1.9 | 50 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 37 | Modeling of magnetic and chemical ordering in binary alloys. <i>Physical Review B</i> , 1984, 30, 1448-1453. | 3.2 | 38 |
| 38 | Experimental and theoretical determination of the metastable Fe-V phase diagram. <i>Physical Review B</i> , 1996, 54, 8958-8961. | 3.2 | 38 |
| 39 | Oxygen order-disorder transition in the superconductor $\text{YBa}_2\text{Cu}_3\text{O}_{6+\delta}$. <i>Physical Review B</i> , 1988, 37, 3678-3680. | 3.2 | 35 |
| 40 | Phase separation and dynamical scaling in borate glasses. <i>Physical Review B</i> , 1986, 34, 2762-2769. | 3.2 | 34 |
| 41 | Phase transitions in Ising square antiferromagnets with first- and second-neighbour interactions. <i>Journal of Physics Condensed Matter</i> , 1994, 6, 9759-9772. | 1.8 | 34 |
| 42 | Electronic-Structure calculation of ordering and segregation energies of transition metal alloys. <i>Acta Metallurgica</i> , 1988, 36, 367-375. | 2.1 | 33 |
| 43 | Mechanical properties of Rh-based L12 intermetallic compounds Rh_3Ti , Rh_3Nb and Rh_3Ta . <i>Intermetallics</i> , 2000, 8, 785-791. | 3.9 | 33 |
| 44 | Magnetic properties and chemical ordering in Co-Pt. <i>Journal of Physics C: Solid State Physics</i> , 1988, 21, L1091-L1096. | 1.5 | 32 |
| 45 | Foundations and Practical Implementations of the Cluster Expansion. <i>Journal of Phase Equilibria and Diffusion</i> , 2017, 38, 238-251. | 1.4 | 32 |
| 46 | Ternary diffusion in multilayer Ag-Au-Cu thin films. <i>Thin Solid Films</i> , 1975, 25, 465-482. | 1.8 | 31 |
| 47 | Application of group theory to the calculation of the configurational entropy in the cluster variation method. <i>Physica A: Statistical Mechanics and Its Applications</i> , 1982, 113, 315-337. | 2.6 | 30 |
| 48 | Calculation of temperature-concentration diagrams by the cv method with Lennard-Jones pair interactions. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 1984, 8, 221-231. | 1.6 | 28 |
| 49 | Characterization of mechanical properties in the Ir-Nb-Zr intermetallic system. <i>Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing</i> , 1993, 170, 169-175. | 5.6 | 28 |
| 50 | Thermal conductivity and thermal expansion of L12 intermetallic compounds based on rhodium. <i>Journal of Alloys and Compounds</i> , 2003, 354, 202-207. | 5.5 | 27 |
| 51 | Reciprocal-space analysis of short-range-order intensities by the cluster-variation method. <i>Physical Review B</i> , 1995, 51, 3429-3437. | 3.2 | 26 |
| 52 | Vibrational free energy in the Ni-Cr system. <i>Computational Materials Science</i> , 1997, 8, 92-99. | 3.0 | 26 |
| 53 | Electronic-structure calculations of binary-alloy phase diagrams. <i>Physical Review B</i> , 1986, 33, 4782-4792. | 3.2 | 25 |
| 54 | Boundary-condition effects on the surface magnetization of Ising ferromagnets. <i>Physical Review B</i> , 1989, 39, 9746-9748. | 3.2 | 25 |

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|----|--|-----|-----------|
| 55 | Thermal conductivity and thermal expansion of Ir ₃ X (X = Ti, Zr, Hf, V, Nb, Ta) compounds for high-temperature applications. <i>Materials Chemistry and Physics</i> , 2003, 80, 385-390. | 4.0 | 25 |
| 56 | Surface first-order phase transitions in fcc Ising ferromagnets. <i>Physical Review Letters</i> , 1987, 58, 1120-1122. | 7.8 | 24 |
| 57 | Experimental determination of pair interaction energies in aCoPt ₃ single crystal and phase-diagram calculations. <i>Physical Review B</i> , 2000, 61, 14975-14983. | 3.2 | 24 |
| 58 | Atomic migration and ordering energies in FePd: Measurement and modeling. <i>Scripta Materialia</i> , 2005, 53, 435-440. | 5.2 | 24 |
| 59 | Diffusion in AgAuPd thin film microcouples. <i>Acta Metallurgica</i> , 1974, 22, 709-719. | 2.1 | 22 |
| 60 | Workshop on thermodynamic modelling of solutions and alloys. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 1997, 21, 219-246. | 1.6 | 22 |
| 61 | General relations between many-body potentials and cluster expansions in multicomponent systems. <i>Journal of Physics Condensed Matter</i> , 2004, 16, 3843-3852. | 1.8 | 21 |
| 62 | Characterization of the deformation behavior of the Cr ₂ Nb ordered intermetallic system. <i>Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing</i> , 1993, 170, 177-183. | 5.6 | 20 |
| 63 | Short range order hardening with second neighbor interactions in fcc solid solutions. <i>Metallurgical and Materials Transactions A - Physical Metallurgy and Materials Science</i> , 1986, 17, 189-194. | 1.4 | 19 |
| 64 | Experimental determination of pair interactions in aFe _{0.804} V _{0.196} single crystal. <i>Physical Review B</i> , 1995, 51, 5760-5767. | 3.2 | 19 |
| 65 | Finite-Size Effects on First-Order Phase Transitions: fcc Binary Alloys. <i>Physical Review Letters</i> , 1986, 57, 360-363. | 7.8 | 18 |
| 66 | Ferromagnetism and chemical ordering in cobalt-iron. <i>Journal of Physics F: Metal Physics</i> , 1988, 18, 767-777. | 1.6 | 18 |
| 67 | On the analysis of diffusion anomalies in Ti alloys. <i>Philosophical Magazine A: Physics of Condensed Matter, Structure, Defects and Mechanical Properties</i> , 1981, 43, 1407-1417. | 0.6 | 17 |
| 68 | Modeling of the Fe-Al phase diagram. <i>Physical Review B</i> , 1988, 38, 11481-11485. | 3.2 | 17 |
| 69 | The effect of the structural energy of Cr on the Ni ⁺ -Cr equilibrium phase diagram. <i>Acta Metallurgica</i> , 1989, 37, 121-127. | 2.1 | 16 |
| 70 | Workshop on thermodynamic models and data for pure elements and other endmembers of solutions. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 1995, 19, 481-498. | 1.6 | 16 |
| 71 | A microscopic theory of binary alloy phase equilibrium. <i>Solid State Communications</i> , 1985, 55, 253-256. | 1.9 | 15 |
| 72 | Modelling of the electrical resistivity of ferromagnetic and paramagnetic intermetallic compounds. <i>Journal of Physics Condensed Matter</i> , 1992, 4, 9053-9066. | 1.8 | 15 |

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|----|---|-----|-----------|
| 73 | The phase diagram of oxygen ordering in YBa ₂ Cu ₃ O _{6+δ} . Physica C: Superconductivity and Its Applications, 1993, 210, 401-407. | 1.2 | 14 |
| 74 | Order-Disorder Transitions under Confinement. Physical Review Letters, 1998, 81, 1146-1149. | 7.8 | 14 |
| 75 | Theoretical study of thermodynamics relevant to tetramethylsilane pyrolysis. Journal of Crystal Growth, 1997, 178, 513-517. | 1.5 | 13 |
| 76 | Stress-Strain Behavior of Nicalon-Fiber-Reinforced Calcium Aluminosilicate Composites under Tensile Fatigue Conditions. Journal of the American Ceramic Society, 1993, 76, 2175-2179. | 3.8 | 12 |
| 77 | Small angle X-ray scattering study of phase separation in glasses using a new position sensitive detector. Nuclear Instruments & Methods in Physics Research, 1983, 208, 489-494. | 0.9 | 11 |
| 78 | Thermodynamic modeling of site occupation in the $\hat{\Gamma}^2$ phase of the Ni–Al–Hf system. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 1989, 108, 159-164. | 5.6 | 10 |
| 79 | Pair correlations and interaction energies in FePd single crystal. European Physical Journal B, 2004, 41, 207-212. | 1.5 | 10 |
| 80 | First principles phase stability study of the Ru–Nb–Zr system. Materials Science & Engineering A: Structural Materials: Properties, Microstructure and Processing, 1993, 170, 161-167. | 5.6 | 8 |
| 81 | Thermodynamic, electronic and magnetic properties of intermetallic compounds through statistical models. Physica Scripta, 1993, T49A, 364-372. | 2.5 | 8 |
| 82 | Simulation of chemical-vapor-deposited silicon carbide for a cold wall vertical reactor. Journal of Crystal Growth, 1997, 178, 505-512. | 1.5 | 8 |
| 83 | Approximate solutions to the cluster variation free energies by the variable basis cluster expansion. Computational Materials Science, 2016, 122, 301-306. | 3.0 | 8 |
| 84 | Modeling of phase diagrams for bcc magnetic alloys. Physical Review B, 1988, 38, 4955-4962. | 3.2 | 7 |
| 85 | Understanding the high-temperature deformation. Metallurgical and Materials Transactions A - Physical Metallurgy and Materials Science, 1992, 23, 3073-3076. | 1.4 | 7 |
| 86 | Ordering in fcc lattices by the Cluster Variation Method and Path Probability Method. Materials Transactions, JIM, 1992, 33, 558-564. | 0.9 | 6 |
| 87 | Thermodynamics of binary alloy thin films. Computational Materials Science, 1997, 8, 79-86. | 3.0 | 6 |
| 88 | Pseudo-three dimensional model of oxygen ordering in YBa ₂ Cu ₃ O _{6+δ} . Solid State Communications, 1991, 79, 151-154. | 1.9 | 5 |
| 89 | Structural energies and phase stability in Ni-Cr alloys. Bulletin of Alloy Phase Diagrams, 1989, 10, 319-326. | 0.2 | 4 |
| 90 | First-principles calculation of the Ni-Cr phase diagram. Solid State Ionics, 1989, 32-33, 714-720. | 2.7 | 4 |

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|-----|---|-----|-----------|
| 91 | Finite-size effects on the phase diagrams of binary alloy films. Solid State Communications, 1998, 107, 285-289. | 1.9 | 4 |
| 92 | Surface ferromagnetism in close-packed structures. Surface Science, 1988, 198, L299-L306. | 1.9 | 3 |
| 93 | Confinement effects in antiferromagnets. Physical Review B, 2000, 62, 1148-1156. | 3.2 | 3 |
| 94 | Cluster-method study of phase separation in He^3 mixtures. Physical Review B, 1986, 33, 5059-5061. | 3.2 | 2 |
| 95 | Continuous sequence of mean-field approximations and critical phenomena. Physica A: Statistical Mechanics and Its Applications, 1994, 209, 257-267. | 2.6 | 2 |
| 96 | Phase Transitions in Confined Antiferromagnets. Physica Status Solidi (B): Basic Research, 2000, 220, 389-394. | 1.5 | 2 |
| 97 | Reply to "Comment on "Cluster expansion and the configurational theory of alloys". Physical Review B, 2017, 95, . | 3.2 | 2 |
| 98 | Investigations of the Co-Pt alloy phase diagram with neutron diffuse scattering, inverse cluster variation method, and Monte Carlo simulations. Physical Review B, 2020, 102, . | 3.2 | 2 |
| 99 | Intermetallic alloys and their composites: The promise and challenges remain. Jom, 1993, 45, 36-36. | 1.9 | 0 |
| 100 | Observation of crack closure in Nicalon/CAS composites. Journal of Materials Science Letters, 1995, 14, 571-573. | 0.5 | 0 |