

# Sergi Vela

## List of Publications by Year in descending order

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papers

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394421  
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477307  
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all docs

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docs citations

61  
times ranked

1293  
citing authors

#	ARTICLE	IF	CITATIONS
1	Heteroatom oxidation controls singlet-triplet energy splitting in singlet fission building blocks. Chemical Communications, 2022, 58, 1338-1341.	4.1	6
2	Donor-Acceptor-Donor Hot Exciton-Triads for High Reverse Intersystem Crossing in OLEDs. Advanced Optical Materials, 2022, 10, .	7.3	7
3	Insights into the magnetism and phase transitions of organic radical-based materials. Journal of Materials Chemistry C, 2021, 9, 10624-10646.	5.5	27
4	Pitfalls on evaluating pair exchange interactions for modelling molecule-based magnetism. Journal of Materials Chemistry C, 2021, 9, 10647-10660.	5.5	7
5	Identifying the Trade-off between Intramolecular Singlet Fission Requirements in Donor-Acceptor Copolymers. Chemistry of Materials, 2021, 33, 2567-2575.	6.7	14
6	Learning the Exciton Properties of Azo-dyes. Journal of Physical Chemistry Letters, 2021, 12, 5957-5962.	4.6	4
7	Tuning the Thermal Stability and Photoisomerization of Azoheteroarenes through Macrocyclic Strain**. Chemistry - A European Journal, 2021, 27, 419-426.	3.3	12
8	Bi-stable spin-crossover in charge-neutral [Fe(R-ptp) <sub>2</sub> ] (ptp =) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 467 Td (2-(1<i>H</i>-p 1022-1031.	3.3	16
9	The Photoisomerization Pathway(s) of Push-Pull Phenylazoheteroarenes**. Chemistry - A European Journal, 2020, 26, 14724-14729.	3.3	6
10	Designing Singlet Fission Candidates from Donor-Acceptor Copolymers. Chemistry of Materials, 2020, 32, 6515-6524.	6.7	27
11	Extremely well isolated two-dimensional spin- $\frac{1}{2} \times \frac{2}{2}$ antiferromagnetic Heisenberg layers with a small exchange coupling in the molecular-based magnet CuPOF. Physical Review B, 2020, 102, .	3.2	8
12	Reversible Magnetic Transition in a Bench-Stable Radical Cation Triggered by Structural Transition in the Magnetically Silent Counteranion. Crystal Growth and Design, 2020, 20, 6296-6301.	3.0	10
13	Structure: function relationships for thermal and light-induced spin-crossover in isomorphous molecular materials. Journal of Materials Chemistry C, 2020, 8, 8420-8429.	5.5	11
14	Assessing Cu <sub>2</sub> L <sub>2</sub> X <sub>4</sub> dimeric moieties as ferromagnetic building blocks in double halide-bridged polymers (X = Cl <sup>-</sup> , Br <sup>-</sup> and L = benzamide). An experimental and computational study. Polyhedron, 2020, 185, 114603.	2.2	2
15	Two different mechanisms of stabilization of regular $\pi$ -stacks of radicals in switchable dithiazolyl-based materials. Journal of Materials Chemistry C, 2020, 8, 5437-5448.	5.5	7
16	Thermal spin crossover in Fe(II) and Fe(III). Accurate spin state energetics at the solid state. Physical Chemistry Chemical Physics, 2020, 22, 4938-4945.	2.8	32
17	Exploring chemical space in the search for improved azoheteroarene-based photoswitches. Physical Chemistry Chemical Physics, 2019, 21, 20782-20790.	2.8	10
18	Deciphering crystal packing effects in the spin crossover of six [Fe <sup>II</sup> (2-pic) <sub>3</sub> ]Cl <sub>2</sub> solvatomorphs. Dalton Transactions, 2019, 48, 1237-1245.	3.3	18

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19	Revising the common understanding of metamagnetism in the molecule-based bisdithiazolyl BDTMe compound. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 12184-12191.	2.8	8
20	Bi-stable spin-crossover characteristics of a highly distorted $[\text{Fe}(\text{1-BPP-COOC})_2(\text{H})_5]_2(\text{ClO}_4)_2 \cdot \text{CH}_3\text{CN}$ complex. <i>Dalton Transactions</i> , 2019, 48, 3825-3830.	2.8	27
21	Controlling the crystallinity and crystalline orientation of "shuttlecock" naphthalocyanine films for near-infrared optoelectronic applications. <i>Journal of Materials Chemistry C</i> , 2018, 6, 1959-1970.	5.5	8
22	Luminescent Dinuclear Copper(I) Complexes as Potential Thermally Activated Delayed Fluorescence (TADF) Emitters: A Theoretical Study. <i>Journal of Physical Chemistry A</i> , 2018, 122, 1413-1421.	2.5	34
23	Pairing-up viologen cations and dications: a microscopic investigation of van der Waals interactions. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 27878-27884.	2.8	10
24	Absorption Spectroscopy and Photophysics of a $\text{Re}^{\text{I}}$ -dppz Probe for DNA-Mediated Charge Transport. <i>Chemistry - A European Journal</i> , 2018, 24, 14425-14435.	3.3	9
25	Cooperativity in Spin Crossover Systems. An Atomistic Perspective on the Devil's Staircase. <i>Inorganic Chemistry</i> , 2018, 57, 9478-9488.	4.0	28
26	A Spin-Crossover Molecular Material Describing Four Distinct Thermal Pathways. <i>Inorganic Chemistry</i> , 2018, 57, 11019-11026.	4.0	19
27	Spin State Chemistry: Modulation of Ligand $\text{p} \rightarrow \text{K} \rightarrow \text{a}$ by Spin State Switching in a $[\text{2} \rightarrow \text{2}]$ Iron(II) Grid-Type Complex. <i>Journal of the American Chemical Society</i> , 2018, 140, 8218-8227.	13.7	63
28	Bistability in Organic Magnetic Materials: A Comparative Study of the Key Differences between Hysteretic and Non-hysteretic Spin Transitions in Dithiazolyl Radicals. <i>Chemistry - A European Journal</i> , 2017, 23, 3479-3489.	3.3	26
29	Electron transport through a spin crossover junction. Perspectives from a wavefunction-based approach. <i>Journal of Chemical Physics</i> , 2017, 146, 064112.	3.0	9
30	Lattice-Solvent Effects in the Spin-Crossover of an Fe(II)-Based Material. The Key Role of Intermolecular Interactions between Solvent Molecules. <i>Inorganic Chemistry</i> , 2017, 56, 4474-4483.	4.0	36
31	Twisting induces ferromagnetism in homometallic clusters. <i>Dalton Transactions</i> , 2017, 46, 11154-11158.	3.3	1
32	A probe of steric ligand substituent effects on the spin crossover of $\text{Fe}(\text{II})$ complexes. <i>Inorganic Chemistry Frontiers</i> , 2017, 4, 1374-1383.	6.0	28
33	Disclosing the Ligand- and Solvent-Induced Changes on the Spin Transition and Optical Properties of Fe(II)-Indazolylpyridine Complexes. <i>Magnetochemistry</i> , 2016, 2, 6.	2.4	10
34	On the zeroth-order hamiltonian for $\text{CASPT}_2$ calculations of spin crossover compounds. <i>Journal of Computational Chemistry</i> , 2016, 37, 947-953.	3.3	36
35	Understanding the Influence of the Electronic Structure on the Crystal Structure of a TTF-PTM Radical Dyad. <i>Journal of Physical Chemistry A</i> , 2016, 120, 10297-10303.	2.5	5
36	Three Redox States of a Diradical Acceptor-Donor-Acceptor Triad: Gating the Magnetic Coupling and the Electron Delocalization. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2234-2239.	4.6	24

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37	Designed intramolecular blocking of the spin crossover of an Fe( $\text{Fe}^{\text{II}}$ ) complex. Dalton Transactions, 2016, 45, 14058-14062.	3.3	15
38	Towards an accurate and computationally-efficient modelling of Fe( $\text{Fe}^{\text{II}}$ )-based spin crossover materials. Physical Chemistry Chemical Physics, 2015, 17, 16306-16314.	2.8	53
39	The origin of the antiferromagnetic behaviour of the charge-transfer compound (HMTTF)[Ni( $\text{mnt}$ ) $_2$ ]. Dalton Transactions, 2015, 44, 608-614.	3.3	5
40	Dynamical effects on the magnetic properties of dithiazolyl bistable materials. Chemical Science, 2015, 6, 2371-2381.	7.4	34
41	Towards the tailored design of benzotriazinyl-based organic radicals displaying a spin transition. Chemical Communications, 2015, 51, 15776-15779.	4.1	16
42	Elucidating the 2D Magnetic Topology of the $\text{Cu}^{\text{II}}$ -Radical $\text{TTTA}^{\text{TM}}$ $\text{Cu}^{\text{II}}$ System. Chemistry - A European Journal, 2014, 20, 7083-7090.	3.3	16
43	On the Importance of Thermal Effects and Crystalline Disorder in the Magnetism of Benzotriazinyl-Derived Organic Radicals. Chemistry - an Asian Journal, 2014, 9, 3612-3622.	3.3	14
44	Insights into the crystal-packing effects on the spin crossover of $[\text{Fe}^{\text{II}}(\text{bpp})]^{2+}$ -based materials. Physical Chemistry Chemical Physics, 2014, 16, 27012-27024.	2.8	57
45	The polymorphism of a triarylphosphine oxide: a case of missing isomers. CrystEngComm, 2014, 16, 8214-8223.	2.6	1
46	The key role of vibrational entropy in the phase transitions of dithiazolyl-based bistable magnetic materials. Nature Communications, 2014, 5, 4411.	12.8	55
47	Linear or Cyclic Clusters of Cu(II) with a Hierarchical Relationship. Inorganic Chemistry, 2014, 53, 3290-3297.	4.0	16
48	A theoretical analysis of the magnetic properties of the low-dimensional copper(II) $\text{X}_2(2\text{-X-3-methylpyridine})_2$ (X = Cl and Br) complexes. Highlights in Theoretical Chemistry, 2014, , 219-230.	0.0	0
49	Dividing the Spoils: Role of Pyrazine Ligands and Perchlorate Counterions in the Magnetic Properties of Bis(pyrazine)diperchloratecopper(II), $[\text{Cu}(\text{pz})_2](\text{ClO}_4)_2$ . Inorganic Chemistry, 2013, 52, 12923-12932.	4.0	22
50	A theoretical analysis of the magnetic properties of the low-dimensional copper(II) $\text{X}_2(2\text{-X-3-methylpyridine})_2$ (X = Cl and Br) complexes. Theoretical Chemistry Accounts, 2013, 132, 1.	1.4	4
51	A theoretical analysis of the magnetic properties of the low dimensional bis(2-chloropyrazine)dichlorocopper(II) molecule-based magnet. Polyhedron, 2013, 64, 163-171.	2.2	2
52	Assigning the dimensionality in low-dimensional materials: A rigorous study of the dimensionality of (2,5-dimethylpyrazine) $\text{CuCl}_2$ . Polyhedron, 2013, 52, 699-705.	2.2	6
53	Tracing the Sources of the Different Magnetic Behavior in the Two Phases of the Bistable (BDTA) $_2$ [Co( $\text{mnt}$ ) $_2$ ] Compound. Inorganic Chemistry, 2012, 51, 8646-8648.	4.0	12
54	A molecular dynamics simulation of methane adsorption in single walled carbon nanotube bundles. Carbon, 2011, 49, 4544-4553.	10.3	46