

# Enric Canadell

## List of Publications by Citations

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L-index

#	Paper	IF	Citations
425	Conceptual aspects of structure-property correlations and electronic instabilities, with applications to low-dimensional transition-metal oxides. <i>Chemical Reviews</i> , <b>1991</b> , 91, 965-1034	68.1	254
424	Analogies between the concepts of molecular chemistry and solid-state physics concerning structural instabilities. Electronic origin of the structural modulations in layered transition metal dichalcogenides. <i>Journal of the American Chemical Society</i> , <b>1992</b> , 114, 9587-9600	16.4	192
423	Hidden fermi surface nesting and charge density wave instability in low-dimensional metals. <i>Science</i> , <b>1991</b> , 252, 96-8	33.3	161
422	Electrical magnetochiral anisotropy in a bulk chiral molecular conductor. <i>Nature Communications</i> , <b>2014</b> , 5, 3757	17.4	132
421	Electronic structure, optical properties, and lattice dynamics in atomically thin indium selenide flakes. <i>Nano Research</i> , <b>2014</b> , 7, 1556-1568	10	132
420	On the band electronic structure of X [M (dmit) <sub>2</sub> ] <sub>2</sub> (X = TTF, (CH <sub>3</sub> ) <sub>4</sub> N ; M = Ni, Pd) molecular conductors and superconductors. <i>Journal De Physique</i> , <b>1989</b> , 50, 2967-2981		132
419	Nanotexturing To Enhance Photoluminescent Response of Atomically Thin Indium Selenide with Highly Tunable Band Gap. <i>Nano Letters</i> , <b>2016</b> , 16, 3221-9	11.5	119
418	Chemical effects on the optical band-gap of heavily doped ZnO:MIII (M=Al,Ga,In): An investigation by means of photoelectron spectroscopy, optical measurements under pressure, and band structure calculations. <i>Physical Review B</i> , <b>2009</b> , 79,	3.3	106
417	Hydrogen-bond tuning of macroscopic transport properties from the neutral molecular component site along the series of metallic organic-inorganic solvates (BEDT-TTF) <sub>4</sub> Re <sub>6</sub> Se <sub>5</sub> Cl <sub>9</sub> .[guest], [guest = DMF, THF, dioxane]. <i>Journal of the American Chemical Society</i> , <b>1993</b> , 115, 4101-4112	16.4	102
416	A Remarkable Family of Rhodium Acetonitrile Compounds Spanning Three Oxidation States and with Nuclearities Ranging from Mononuclear and Dinuclear to One-Dimensional Chains. <i>Journal of the American Chemical Society</i> , <b>1999</b> , 121, 8005-8016	16.4	100
415	Computing the Properties of Materials from First Principles with SIESTA. <i>Structure and Bonding</i> , <b>2004</b> , 103-170	0.9	96
414	A single-component molecular metal based on a thiazole dithiolate gold complex. <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 16961-7	16.4	92
413	Electronic structure of transition-metal borides with the AlB <sub>2</sub> structure. <i>Journal of the American Chemical Society</i> , <b>1986</b> , 108, 6561-6568	16.4	89
412	Chiral molecular metals: syntheses, structures, and properties of the AsF <sub>6</sub> <sup>-</sup> salts of racemic (+/-)-, (R)-, and (S)-tetrathiafulvalene-oxazoline derivatives. <i>Journal of the American Chemical Society</i> , <b>2005</b> , 127, 5748-9	16.4	86
411	Importance of short interlayer Te-Te contacts for the structural distortions and physical properties of CdI <sub>2</sub> -type layered transition-metal ditellurides. <i>Journal of Solid State Chemistry</i> , <b>1992</b> , 99, 189-199	3.3	86
410	Single crystalline commensurate metallic assemblages of pi-slabs and CdI <sub>2</sub> -type layers: synthesis and properties of beta-(EDT-TTF-I <sub>2</sub> ) <sub>2</sub> [Pb <sub>5/6</sub> square 1/6I <sub>2</sub> ](3) and beta-(EDT-TTF-I <sub>2</sub> ) <sub>2</sub> [Pb <sub>2/3+x</sub> Ag <sub>1/3-2x</sub> square xI <sub>2</sub> ] <sub>3</sub> , x = 0.05. <i>Journal of the American Chemical Society</i> , <b>2003</b> , 125, 3295-301	16.4	85
409	Hal-Hal interactions in a series of three isostructural salts of halogenated tetrathiafulvalenes. Contribution of the halogen atoms to the HOMO-HOMO overlap interactions. <i>Journal of Materials Chemistry</i> , <b>2001</b> , 11, 1570-1575		84

408	Single-component magnetic conductors based on Mo <sub>3</sub> S <sub>7</sub> trinuclear clusters with outer dithiolate ligands. <i>Journal of the American Chemical Society</i> , <b>2004</b> , 126, 12076-83	16.4	83
407	Unprecedented Conversion of a Compound with Metal-Metal Bonding into a Solvated Molecular Wire. <i>Angewandte Chemie International Edition in English</i> , <b>1996</b> , 35, 2772-2774		82
406	Concerning the band structure of D(M(dmit) <sub>2</sub> ) <sub>2</sub> (D=TTF,Cs,NMe <sub>4</sub> ); M=Ni,Pd) molecular conductors and superconductors: Role of the M(dmit) <sub>2</sub> Homo and LUMO. <i>Solid State Communications</i> , <b>1990</b> , 75, 633-638	16.4	82
405	Crystalline arrays of pairs of molecular rotors: correlated motion, rotational barriers, and space-inversion symmetry breaking due to conformational mutations. <i>Journal of the American Chemical Society</i> , <b>2013</b> , 135, 9366-76	16.4	81
404	CaFeO <sub>2</sub> : a new type of layered structure with iron in a distorted square planar coordination. <i>Journal of the American Chemical Society</i> , <b>2009</b> , 131, 221-9	16.4	80
403	(EDT-TTF-CONH <sub>2</sub> ) <sub>6</sub> [Re <sub>6</sub> Se <sub>8</sub> (CN) <sub>6</sub> ], a metallic Kagome-type organic-inorganic hybrid compound: electronic instability, molecular motion, and charge localization. <i>Journal of the American Chemical Society</i> , <b>2005</b> , 127, 11785-97	16.4	78
402	Molecular Metals Based on BEDT-TTF Radical Cation Salts with Magnetic Metal Oxalates as Counterions: $\text{A}(\text{BEDT-TTF})_4\text{A}[\text{M}(\text{C}_2\text{O}_4)_3]\cdot\text{DMF}$ (A = NH <sub>4</sub> <sup>+</sup> , K <sup>+</sup> ; M = Cr <sup>III</sup> , Fe <sup>III</sup> ). <i>Advanced Functional Materials</i> , <b>2003</b> , 13, 403-411	15.6	77
401	Directing the Structures and Collective Electronic Properties of Organic Conductors: The Interplay of $\pi$ - $\pi$ Overlap Interactions and Hydrogen Bonds. <i>Chemistry - A European Journal</i> , <b>1999</b> , 5, 2971-2976	4.8	77
400	An Efficient, Redox-Enhanced Pair of Hydrogen-Bond Tweezers for Chloride Anion Recognition, a Key Synthon in the Construction of a Novel Type of Organic Metal based on the Secondary Amide-Functionalized Ethylenedithiotetrathiafulvalene, $\text{A}(\text{EDT-TTF-CONHMe})_2[\text{Cl}\cdot\text{H}_2\text{O}]$ . <i>Chemistry of Materials</i> , <b>2000</b> , 12, 1898-1904	9.6	71
399	Regioselectivity of radical attacks on substituted olefins. Application of the state-correlation-diagram (SCD) model. <i>Journal of the American Chemical Society</i> , <b>1990</b> , 112, 1446-1452	16.4	71
398	Chirality driven metallic versus semiconducting behavior in a complete series of radical cation salts based on dimethyl-ethylenedithio-tetrathiafulvalene (DM-EDT-TTF). <i>Journal of the American Chemical Society</i> , <b>2013</b> , 135, 17176-86	16.4	70
397	Theoretical study of the electrical behavior of one-dimensional metallophthalocyanines and related metallomacrocyclic compounds. <i>Inorganic Chemistry</i> , <b>1984</b> , 23, 573-579	5.1	70
396	Anisotropic chemical pressure effects in single-component molecular metals based on radical dithiolene and diselenolene gold complexes. <i>Journal of the American Chemical Society</i> , <b>2012</b> , 134, 17138-48	16.4	65
395	The [(DT-TTF) <sub>2</sub> M(mnt) <sub>2</sub> ] Family of Radical Ion Salts: From a Spin Ladder to Delocalised Conduction Electrons That Interact with Localised Magnetic Moments. <i>Chemistry - A European Journal</i> , <b>1999</b> , 5, 2025-2039	4.8	65
394	Experimental and theoretical study of band structure of InSe and In <sub>1-x</sub> GaxSe (x. <i>Physical Review B</i> , <b>2001</b> , 63,	3.3	64
393	Structural and electronic properties of SrFeO <sub>2</sub> from first principles. <i>Physical Review B</i> , <b>2008</b> , 78,	3.3	63
392	Theoretical analysis of bonding in monomeric and polymeric C <sub>5</sub> H <sub>5</sub> M compounds. <i>Organometallics</i> , <b>1984</b> , 3, 759-764	3.8	63
391	Extended polymetallic sandwich compounds. <i>Organometallics</i> , <b>1985</b> , 4, 805-815	3.8	63

- 390 Essential role of anions in the charge ordering transition of  $\text{E}(\text{BEDT-TTF})_2\text{I}_3$ . *Physical Review B*, **2012**, 85, 3-3 62
- 389 Electronic structure of transition metal complex-based molecular metals and superconductors. *Coordination Chemistry Reviews*, **1999**, 185-186, 629-651 23.2 60
- 388 A chiral molecular conductor: synthesis, structure, and physical properties of  $[\text{ET}]_3[\text{Sb}_2(\text{L-tart})_2]\cdot\text{CH}_3\text{CN}$  (ET = bis(ethylenedithio)tetrathiafulvalene; L-tart = (2R,3R)-(+)-tartrate). *Inorganic Chemistry*, **2004**, 43, 8072-7 5.1 59
- 387 Singular crystalline beta'-layered topologies directed by ribbons of self-complementary amide...amide ring motifs in  $[\text{EDT-TTF}(\text{CONH}(2))(2)](2)\text{X}$  (X =  $\text{HSO}_4(-)$ ,  $\text{ClO}_4(-)$ ,  $\text{ReO}_4(-)$ ,  $\text{AsF}_6(-)$ ): coupled activation of ribbon curvature, electron interactions, and magnetic susceptibility. *Journal of the American Chemical Society*, **2003**, 125, 11583-90 16.4 59
- 386 Band electronic structure study of the electronic instability in the Magneli phase molybdenum oxide  $[\text{Mo}_4\text{O}_{11}]$ . *Inorganic Chemistry*, **1989**, 28, 1466-1472 5.1 59
- 385 O?S vs. N?S intramolecular nonbonded interactions in neutral and radical cation salts of TTF-oxazoline derivatives: synthesis, theoretical investigations, crystalline structures, and physical properties. *New Journal of Chemistry*, **2007**, 31, 1468 3.6 56
- 384 First-principles study of the neutral molecular metal  $\text{Ni}(\text{tmdt})_2$ . *Physical Review B*, **2002**, 65, 3-3 55
- 383 Crystal symmetry and pressure effects on the valence band structure of  $\text{E}(\text{InSe})$  and  $\text{E}(\text{GaSe})$ : Transport measurements and electronic structure calculations. *Physical Review B*, **2005**, 71, 3-3 54
- 382 A construction principle of the  $\text{E}(\text{phase})$  based on the efficient (O?H)donor  $\text{D}(\text{anion})$  structural functionality: The examples of  $\text{E}(\text{EDT-TTF}(\text{CH}_2\text{OH}))_2\text{X}$  (X = ClO and ReO). *Advanced Materials*, **1992**, 4, 579-581 24 54
- 381 Band electronic structure of the lithium molybdenum purple bronze  $\text{Li}_{0.9}\text{Mo}_6\text{O}_{17}$ . *Journal of the American Chemical Society*, **1988**, 110, 358-363 16.4 54
- 380 Symmetry control of the coloring problem: the electronic structure of  $\text{MB}_2\text{C}_2$  (M = calcium, lanthanum, ...). *Journal of the American Chemical Society*, **1986**, 108, 3971-3976 16.4 54
- 379 Fluorine Segregation Controls the Solid-State Organization and Electronic Properties of Ni and Au Dithiolene Complexes: Stabilization of a Conducting Single-Component Gold Dithiolene Complex. *Advanced Functional Materials*, **2002**, 12, 693-698 15.6 53
- 378 A Genuine Quarter-Filled Band Mott Insulator,  $(\text{EDT-TTF-CONMe}_2)_2\text{AsF}_6$ : Where the Chemistry and Physics of Weak Intermolecular Interactions Act in Unison. *Advanced Materials*, **2003**, 15, 1251-1254 24 51
- 377 Complete series of chiral paramagnetic molecular conductors based on tetramethyl-bis(ethylenedithio)-tetrathiafulvalene (TM-BEDT-TTF) and Chloranilate-bridged heterobimetallic honeycomb layers. *Inorganic Chemistry*, **2015**, 54, 3643-53 5.1 50
- 376 Characterization of the unoccupied and partially occupied states of TTF-TCNQ by XANES and first-principles calculations. *Physical Review B*, **2003**, 68, 3-3 50
- 375 Nature of holes, oxidation states, and hypervalency in covellite ( $\text{CuS}$ ). *Inorganic Chemistry*, **2014**, 53, 12402-6 9.6 49
- 374 Design and evaluation of a crystalline hybrid of molecular conductors and molecular rotors. *Journal of the American Chemical Society*, **2012**, 134, 7880-91 16.4 48
- 373 An Organic Spin-Ladder Molecular Material. *Angewandte Chemie International Edition in English*, **1997**, 36, 2324-2326 48

372	Comparison of the electronic structures of layered transition-metal trichalcogenides thallium triselenide, thallium trisulfide and niobium triselenide. <i>Inorganic Chemistry</i> , <b>1990</b> , 29, 1401-1407	5.1	47
371	Charge-density-wave instabilities expected in monophosphate tungsten bronzes. <i>Physical Review B</i> , <b>1991</b> , 43, 1894-1902	3.3	47
370	Structural isomerism in crystals of redox-active secondary ortho-diamides: the role of competing intra- and intermolecular hydrogen bonds in directing crystalline topologies. <i>Chemistry - A European Journal</i> , <b>2004</b> , 10, 4498-511	4.8	45
369	Fused and linked deltahedral clusters in the chemistry of the Group 13 elements. <i>Journal of the American Chemical Society</i> , <b>1990</b> , 112, 7207-7217	16.4	45
368	Charge ordering, symmetry and electronic structure issues and Wigner crystal structure of the quarter-filled band Mott insulators and high pressure metals $\chi(\text{EDT-TTF-CONMe}_2)_2\text{X}$ , X = Br and AsF <sub>6</sub> . <i>Journal of Materials Chemistry</i> , <b>2009</b> , 19, 6980		44
367	Order versus disorder in chiral tetrathiafulvalene-oxazoline radical-cation salts: structural and theoretical investigations and physical properties. <i>Chemistry - A European Journal</i> , <b>2010</b> , 16, 528-37	4.8	44
366	Superstructures of donor packing arrangements in a series of molecular charge transfer salts. <i>Chemical Communications</i> , <b>2004</b> , 18-9	5.8	44
365	Antiperovskite Structure with Ternary Tetrathiafulvalenium Salts: Construction, Distortion, and Antiferromagnetic Ordering. <i>Angewandte Chemie International Edition in English</i> , <b>1991</b> , 30, 1498-1500		43
364	Electronic instabilities of the quasi-two-dimensional monophosphate tungsten bronze P <sub>4</sub> W <sub>12</sub> O <sub>44</sub> . <i>Physical Review B</i> , <b>1989</b> , 39, 12969-12972	3.3	43
363	Band electronic structure of the purple potassium molybdenum bronze K <sub>0.9</sub> Mo <sub>6</sub> O <sub>17</sub> . <i>Journal of the American Chemical Society</i> , <b>1987</b> , 109, 6308-6313	16.4	43
362	Global instability index optimizations for the localization of mobile protons. <i>Solid State Ionics</i> , <b>2004</b> , 168, 281-290	3.3	42
361	Solution Chemistry of Chalcohalide Hexanuclear Rhenium Cluster Monoanions: Substitution Reactions and Structural and LSIMS Characterization of the Heterosubstituted Cluster Dianions, (n-Bu <sub>4</sub> N) <sub>2</sub> [Re <sub>6</sub> Q <sub>5</sub> ECI <sub>8</sub> ] (Q = S, E = O, S, Se; Q = Se, E = S, Se, Te). <i>Inorganic Chemistry</i> , <b>1995</b> , 34, 5307-5313	5.1	42
360	Superconductivity and magnetic field induced spin density waves in the (TMTTF) <sub>2</sub> X family. <i>Journal De Physique, I</i> , <b>1994</b> , 4, 1539-1549		42
359	Concerning the Structure of Hydrogen Molybdenum Bronze Phase III. A Combined Theoretical/Experimental Study. <i>Chemistry of Materials</i> , <b>2005</b> , 17, 5957-5969	9.6	41
358	Electronic origin of the thermochromic effect in 2,2',5,5'-tetramethylbistibole. <i>Journal of the American Chemical Society</i> , <b>1982</b> , 104, 3876-3879	16.4	41
357	Crystal Structure of the Host Lattices of the Superconductors Li <sub>x</sub> MNX (M = Zr, Hf; X = Cl, Br). <i>Chemistry of Materials</i> , <b>1999</b> , 11, 203-206	9.6	40
356	Structural and electronic properties of Cs(Pd(dmit) <sub>2</sub> ) <sub>2</sub> . <i>Journal of Physics Condensed Matter</i> , <b>1991</b> , 3, 933-954	1.8	40
355	Conducting Anilate-Based Mixed-Valence Fe(II)Fe(III) Coordination Polymer: Small-Polaron Hopping Model for Oxalate-Type Fe(II)Fe(III) 2D Networks. <i>Journal of the American Chemical Society</i> , <b>2018</b> , 140, 12611-12621	16.4	40

- 354 On the electronic structure of MPS3 phases. *Inorganic Chemistry*, **1987**, 26, 963-965 5.1 39
- 353 Hybrid molecular materials based upon organic pi-electron donors and metal complexes. Radical salts of bis(ethylenethia)tetrathiafulvalene (BET-TTF) with the octahedral anions hexacyanoferrate(III) and nitroprusside. The first kappa phase in the BET-TTF family. *Inorganic Chemistry*, **2001**, 40, 2521-2528 5.1 38
- 352 Synthesis, Crystal Structure, Electrical Properties and Electronic Band Structure of (NH<sub>4</sub>Me<sub>4</sub>-y)<sub>x</sub>[M(dmit)<sub>2</sub>] Complexes (M = Ni, Pd, Pt; dmit<sup>2-</sup> = 2-Thioxo-1,3-dithiole-4,5-dithiolato). *Inorganic Chemistry*, **1994**, 33, 3401-3414 5.1 38
- 351 Role of p-d and s-d interactions in the electronic structure and band gap of Zn<sub>1-x</sub>M<sub>x</sub>O (M=Cr, Mn, Fe, Co, Ni, and Cu): Photoelectron and optical spectroscopy and first-principles band structure calculations. *Physical Review B*, **2012**, 86, 3.3 37
- 350 Bis- and tetrakis(1,4-dithiafulven-6-yl)-substituted tetrathiafulvalenes and dihydrotetrathiafulvalenes: a novel class of planar donor molecules with multiple redox functionalities and the demonstration of a novel type of two-dimensional association in the solid state. *Chemistry of Materials*, **1993**, 5, 1196-1198 9.6 37
- 349 Localization versus Delocalization in Chiral Single Component Conductors of Gold Bis(dithiolene) Complexes. *Journal of the American Chemical Society*, **2016**, 138, 6838-51 16.4 37
- 348 Structural diversity and physical properties of paramagnetic molecular conductors based on bis(ethylenedithio)tetrathiafulvalene (BEDT-TTF) and the tris(chloranilato)ferrate(III) complex. *Inorganic Chemistry*, **2014**, 53, 7028-39 5.1 36
- 347 First-principles study of the blue bronze K<sub>0.3</sub>MoO<sub>3</sub>. *Physical Review B*, **2002**, 65, 3.3 36
- 346 Metal Complexes of Dithiolate Ligands: 5,6-Dihydro-1,4-dithiin-2,3-dithiolato (dddt(2-)), 5,7-Dihydro-1,4,6-trithiin-2,3-dithiolato (dtdt(2-)), and 2-Thioxo-1,3-dithiole-4,5-dithiolato (dmit(2-)). Synthesis, Electrochemical Studies, Crystal and Electronic Structures, and Conducting Properties. *Inorganic Chemistry*, **1996**, 35, 3856-3873 5.1 36
- 345 STRUCTURAL AND ELECTRONIC ORIGIN OF THE HIDDEN NESTING AND CHARGE DENSITY WAVES IN TRANSITION METAL OXIDES AND BRONZES. *International Journal of Modern Physics B*, **1993**, 07, 4005-4043 1.1 36
- 344 Bond cleavage of the solvated methyl chloride anion: a primary electrochemical event. *Journal of the American Chemical Society*, **1980**, 102, 855-857 16.4 36
- 343 Stable Metallic State of a Neutral-Radical Single-Component Conductor at Ambient Pressure. *Journal of the American Chemical Society*, **2018**, 140, 6998-7004 16.4 35
- 342 A neutral zwitterionic molecular solid. *Chemistry - A European Journal*, **2010**, 16, 14051-9 4.8 35
- 341 Antifluorite-type lithium chromium oxide nitrides: synthesis, structure, order, and electrochemical properties. *Inorganic Chemistry*, **2004**, 43, 7050-60 5.1 35
- 340 Activation of C-H... Halogen (Cl, Br, and I) hydrogen bonds at the organic/inorganic interface in fluorinated tetrathiafulvalenes salts. *Chemistry - A European Journal*, **2001**, 7, 2635-43 4.8 34
- 339 Angle-resolved photoemission study and first-principles calculation of the electronic structure of GaTe. *Physical Review B*, **2002**, 65, 3.3 34
- 338 Band electronic structure study of the semimetallic properties and the anisotropic resistivity hump in zirconium tritelluride. *Journal of the American Chemical Society*, **1988**, 110, 104-108 16.4 34
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336	Electronic structure of 2 H -NbSe 2 single-layers in the CDW state. <i>2D Materials</i> , <b>2016</b> , 3, 035028	5.9	33
335	Donor-π interactions in quarter-filled low-dimensional organic conductors. <i>Materials Horizons</i> , <b>2018</b> , 5, 590-640	14.4	33
334	The 8 : 1 : 1 ternary hybrid framework in the system [EDT-TTF +][1,4-bis(iodoethynyl)benzene][Re6Se8(CN)6]4-: dual noncovalent expression of the octahedral halogen-bond hexa-acceptor nanonode. <i>Chemical Communications</i> , <b>2008</b> , 2194-6	5.8	33
333	Quasiparticle spectra of 2H-NbSe2: Two-band superconductivity and the role of tunneling selectivity. <i>Physical Review B</i> , <b>2015</b> , 92,	3.3	32
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331	(EDT-TTF-I2)2PbI3·H2O: an ambient pressure metal with a π donor slab topology. <i>Journal of Materials Chemistry</i> , <b>2004</b> , 14, 135-137		32
330	Donor Slab Robustness and Band Filling Variations in BDT-TP-based Molecular Conductors: π(BDT-TP)6[Re6S6Cl8]·(CH2Cl)2 and π(BDT-TP)6 [Mo6Cl14]·(CH2Cl)2. <i>Advanced Materials</i> , <b>2000</b> , 12, 436-439	24	32
329	Origin of metal clustering in transition-metal chalcogenide layers MX2 (M = Nb, Ta, Mo, Re; X = S, Se). <i>Journal of the American Chemical Society</i> , <b>1989</b> , 111, 3778-3782	16.4	32
328	Role of large but defective deltahedra in the structural chemistry of very complex solid borides and gallides. <i>Inorganic Chemistry</i> , <b>1991</b> , 30, 1991-1998	5.1	32
327	Metallic versus nonmetallic properties of ternary chalcogenides: tantalum metal selenide, Ta2MSe7 (M = nickel, platinum), and tantalum nickel chalcogenide, Ta2NiX5 (X = sulfide, selenide). <i>Inorganic Chemistry</i> , <b>1987</b> , 26, 3974-3976	5.1	32
326	High-pressure electronic structure and phase transitions in monoclinic InSe: X-ray diffraction, Raman spectroscopy, and density functional theory. <i>Physical Review B</i> , <b>2008</b> , 77,	3.3	31
325	Intramolecular mixed-valence state through silicon or germanium double bridges in rigid bis(tetrathiafulvalenes). <i>Chemistry - A European Journal</i> , <b>2007</b> , 13, 5394-400	4.8	31
324	A chirality-induced alpha phase and a novel molecular magnetic metal in the BEDT-TTF/tris(croconate)ferrate(III) hybrid molecular system. <i>Chemical Communications</i> , <b>2006</b> , 4931-3	5.8	31
323	Reversible Control of Crystalline Rotors by Squeezing Their Hydrogen Bond Cloud Across a Halogen Bond-Mediated Phase Transition. <i>Crystal Growth and Design</i> , <b>2014</b> , 14, 3375-3383	3.5	30
322	Magnetic Properties and New Structural Classification of Molybdenum Phosphates Containing Mo(V). <i>Chemistry of Materials</i> , <b>1997</b> , 9, 68-75	9.6	30
321	New BEDT-TTF/[Fe(C5O5)3]3- hybrid system: synthesis, crystal structure, and physical properties of a chirality-induced alpha phase and a novel magnetic molecular metal. <i>Inorganic Chemistry</i> , <b>2007</b> , 46, 4446-57	5.1	30
320	Nature of the charge density wave images of layered tantalum dichalcogenides 1T-TaX2 (X = sulfur, selenium) in scanning tunneling and atomic force microscopy. <i>Journal of the American Chemical Society</i> , <b>1993</b> , 115, 3760-3765	16.4	30
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- 318 Synthesis, characterization, and electronic structure of Ba<sub>5</sub>In<sub>4</sub>Bi<sub>5</sub>: an acentric and one-electron deficient phase. *Chemistry - A European Journal*, **2004**, 10, 3615-21 4.8 29
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