

# Lijie Li

## List of Publications by Citations

**Source:** <https://exaly.com/author-pdf/7555184/lijie-li-publications-by-citations.pdf>

**Version:** 2024-04-17

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

63

papers

567

citations

15

h-index

20

g-index

69

ext. papers

800

ext. citations

3.5

avg, IF

4.11

L-index

#	Paper	IF	Citations
63	Effects of Additives on HNIW Crystal Morphology and Impact Sensitivity. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2012</b> , 37, 77-82	1.7	44
62	Preparation, characterization and thermal risk evaluation of dihydroxylammonium 5, 5'-bistetrazole-1, 1'-diolate based polymer bonded explosive. <i>Journal of Hazardous Materials</i> , <b>2017</b> , 338, 208-217	12.8	36
61	Thermolysis, nonisothermal decomposition kinetics, calculated detonation velocity and safety assessment of dihydroxylammonium 5, 5'-bistetrazole-1, 1'-diolate. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2016</b> , 126, 473-480	4.1	28
60	Nitrogen-rich 4,4'-azo bis(1,2,4-triazolone) salts--the synthesis and promising properties of a new family of high-density insensitive materials. <i>Dalton Transactions</i> , <b>2016</b> , 45, 3590-8	4.3	24
59	Polytetrafluoroethylene-assisted N/F co-doped hierarchically porous carbon as a high performance electrode for supercapacitors. <i>Journal of Colloid and Interface Science</i> , <b>2019</b> , 545, 25-34	9.3	22
58	Surface Modification for Improving the Photocatalytic Polymerization of 3,4-Ethylenedioxythiophene over Inorganic Lead Halide Perovskite Quantum Dots. <i>ACS Applied Materials &amp; Interfaces</i> , <b>2020</b> , 12, 451-460	9.5	21
57	Morphology control of 3-nitro-1,2,4-triazole-5-one (NTO) by molecular dynamics simulation. <i>CrystEngComm</i> , <b>2018</b> , 20, 6252-6260	3.3	20
56	Prediction of crystal morphology of 3,4-Dinitro-1H-pyrazole (DNP) in different solvents. <i>Journal of Molecular Graphics and Modelling</i> , <b>2017</b> , 75, 62-70	2.8	19
55	Thermal hazard assessment of 4,10-dinitro-2,6,8,12-tetraoxa-4,10-diazaisowurtzitane (TEX) by accelerating rate calorimeter (ARC). <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2016</b> , 126, 467-471	4.1	19
54	Polymer-based lightweight materials for electromagnetic interference shielding: a review. <i>Journal of Materials Science</i> , <b>2021</b> , 56, 6549-6580	4.3	19
53	Molecular dynamic simulations on TKX-50/HMX cocrystal. <i>RSC Advances</i> , <b>2017</b> , 7, 6795-6799	3.7	18
52	Quantitative Determination of H phase in polymorphic HNIW using X-ray Diffraction Patterns. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2008</b> , 33, 467-471	1.7	17
51	A simple ratiometric and colorimetric chemosensor for the selective detection of fluoride in DMSO buffered solution. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , <b>2016</b> , 153, 194-198	4.4	15
50	Squaramide-based lab-on-a-molecule for the detection of silver ion and nitroaromatic explosives. <i>RSC Advances</i> , <b>2015</b> , 5, 96665-96669	3.7	15
49	The effect of solution conditions on the crystal morphology of HMX by molecular dynamics simulations. <i>Journal of Crystal Growth</i> , <b>2019</b> , 507, 38-45	1.6	15
48	Molecular dynamics simulations on miscibility, glass transition temperature and mechanical properties of PMMA/DBP binary system. <i>Journal of Molecular Graphics and Modelling</i> , <b>2018</b> , 84, 182-188	2.8	13
47	Preparation, nonisothermal decomposition kinetics, heat capacity, and safety parameters of TKX-50-based PBX. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2018</b> , 131, 3193-3199	4.1	13

46	The crossover behavior of bromine species in the metal-free flow battery. <i>Journal of Applied Electrochemistry</i> , <b>2017</b> , 47, 261-272	2.6	12
45	Preparation and Characterization of Cyclotrimethylenetrinitramine (RDX) with Reduced Sensitivity. <i>Materials</i> , <b>2017</b> , 10,	3.5	12
44	The study of external growth environments on the crystal morphology of HNIW by molecular dynamics simulation. <i>Journal of Materials Science</i> , <b>2018</b> , 53, 12921-12936	4.3	11
43	First-principles calculation of quantum capacitance of metals doped graphenes and nitrogen/metals co-doped graphenes: designing strategies for supercapacitor electrodes. <i>Journal of Materials Science</i> , <b>2019</b> , 54, 483-492	4.3	11
42	Thermal behavior and thermo-kinetic studies of 5,5?-bistetrazole-1,1?-diolate (1,1-BTO). <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2017</b> , 129, 1265-1270	4.1	10
41	Dissolution Properties of Dihydroxylammonium 5,5?-Bistetrazole-1,1?-diolate and Disodium 5,5?-Bistetrazole-1,1?-diolate in Water. <i>Journal of Energetic Materials</i> , <b>2016</b> , 34, 416-425	1.6	10
40	Preparation, Crystal Structure and Properties of a New Crystal Form of Diammonium 5,5?-bistetrazole-1,1?-diolate. <i>Chinese Journal of Chemistry</i> , <b>2015</b> , 33, 1229-1234	4.9	10
39	Preparation and performances characterization of HNIW/NTO-based high-energetic low vulnerable polymer-bonded explosive. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2020</b> , 139, 3589-3602	4.1	10
38	Thermal decomposition mechanism study of 3-nitro-1,2,4-triazol-5-one (NTO): Combined TG-FTIR-MS techniques and ReaxFF reactive molecular dynamics simulations. <i>Fuel</i> , <b>2021</b> , 295, 120655	7.1	10
37	Study on a novel high energetic and insensitive munitions formulation: TKX-50 based melt cast high explosive. <i>RSC Advances</i> , <b>2017</b> , 7, 31485-31492	3.7	9
36	The novel compound dimethylamine-5,5?-bistetrazole-1,1?-diolate: crystal structure, thermal investigation, safety evaluation and theoretical studies. <i>RSC Advances</i> , <b>2017</b> , 7, 18523-18528	3.7	9
35	Effects of the N/S codoping configuration and ternary doping on the quantum capacitance of graphene. <i>Journal of Materials Science</i> , <b>2019</b> , 54, 8995-9003	4.3	8
34	Reactive molecular dynamics simulations on the thermal decompositions and oxidations of TKX-50 and twinned TKX-50. <i>CrystEngComm</i> , <b>2020</b> , 22, 2593-2600	3.3	7
33	Theoretical study of the heats of formation, detonation properties, and bond dissociation energies of substituted bis-1,2,4-triazole compounds. <i>Journal of Molecular Modeling</i> , <b>2018</b> , 24, 85	2	6
32	Preparation and thermal properties study of HNIW/FOX-7 based high energy polymer bonded explosive (PBX) with low vulnerability to thermal stimulations. <i>Journal of Energetic Materials</i> , <b>2020</b> , 38, 83-97	1.6	6
31	Thermal decomposition and thermal stability of potassium 3,3?-dinitrimino-5,5?-bis(1H-1,2,4-triazole). <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2018</b> , 133, 1563-1569	4.1	5
30	The Critical Role of Oxygen-Containing Functional Groups in the Etching Behavior of Activators to Carbon Materials. <i>ACS Sustainable Chemistry and Engineering</i> , <b>2021</b> , 9, 1646-1655	8.3	5
29	Dissolution thermodynamics of dihydroxylammonium 5,5?-bistetrazole-1,1?-diolate in water at T = (298.15, 303.15, 308.15 and 313.15 K). <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2017</b> , 128, 1875-1880	4.1	4

28	Density Functional Theory (DFT) Study on the Structures and Energetic Properties of Isomers of Tetranitro-bis-1,2,4-triazoles. <i>ACS Omega</i> , <b>2020</b> , 5, 19464-19468	3.9	4
27	Investigation into the Temperature Adaptability of HNIW-based PBXs. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2019</b> , 44, 327-336	1.7	4
26	Molecular dynamic simulations for FOX-7 and FOX-7 based PBXs. <i>Journal of Molecular Modeling</i> , <b>2018</b> , 24, 145	2	4
25	Preparation, crystal structure, thermal behavior, and theoretical studies of N,N'-dinitro-4,4'-azo-bis(1,2,4-triazolone) (DNZTO). <i>Zeitschrift Fur Naturforschung - Section B Journal of Chemical Sciences</i> , <b>2016</b> , 71, 197-204	1	3
24	Miscibility, Glass Transition Temperature and Mechanical Properties of NC/DBP Binary Systems by Molecular Dynamics. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2018</b> , 43, 559-567	1.7	3
23	Measurement and Correlation of Solubilities of 5,5'-Dinitramino-3,3'-bi[1,2,4-triazolate] Carbohydrazide Salt (CBNT) in Various Pure Solvents and a Binary Mixture (Dimethyl Sulfoxide + Water) from 298.15 to 343.15 K. <i>Journal of Chemical &amp; Engineering Data</i> , <b>2019</b> , 64, 3874-3881	2.8	3
22	Heat effects of NTO synthesis in nitric acid solution. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2017</b> , 128, 301-310	4.1	3
21	Synthesis and Characterization of 1,5-Dinitro-2,6-bis(trinitromethyl)-3a,4a,7a,8a-tetrahydro-[1,4]dioxino[2,3-d:5,6-d']diimidazole (DNTDI). <i>Propellants, Explosives, Pyrotechnics</i> , <b>2013</b> , 38, 658-664	1.7	3
20	A molecular dynamics study and detonation parameters calculation of 5,5'-dinitramino-3,3'-bi[1,2,4-triazolate] carbohydrazide salt (CBNT) and its PBXs. <i>Journal of Energetic Materials</i> , <b>2020</b> , 38, 283-294	1.6	3
19	Solubilities of 2,6-Diamino-3,5-dinitropyrazine-1-oxide in the Binary Mixtures of DMSO+H <sub>2</sub> O, DMF+H <sub>2</sub> O and NMP+H <sub>2</sub> O in the Temperature Range from 293.15 to 323.15 K under the Atmospheric Pressure. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2020</b> , 45, 503-508	1.7	3
18	Empirical Kinetics Equation of the Synthesis of NTO in Nitric Acid. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2016</b> , 41, 1085-1091	1.7	3
17	Thermal safety assessment and thermo-kinetic parameters of 5,5'-dinitramino-3,3'-bi[1,2,4-triazolate] carbohydrazide salt (CBNT). <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2021</b> , 144, 647-655	4.1	3
16	Dissolution properties of 5,5'-bistetrazole-1, 1'-dihydroxy and disodium 5,5'-bistetrazole-1, 1'-diolate in dimethyl sulfoxide. <i>Journal of Thermal Analysis and Calorimetry</i> , <b>2017</b> , 128, 615-620	4.1	2
15	Theoretical study on the weak interaction and energy performance of nitroformate salts and nitroformate-based propellant formulations. <i>Journal of Molecular Modeling</i> , <b>2019</b> , 25, 285	2	2
14	Study on Cellulose Acetate Butyrate/Plasticizer Systems by Molecular Dynamics Simulation and Experimental Characterization. <i>Polymers</i> , <b>2020</b> , 12,	4.5	2
13	The crystal structure and thermal analysis of ZTO and its solvent adducts. <i>Research on Chemical Intermediates</i> , <b>2016</b> , 42, 4333-4340	2.8	2
12	Ameliorating the performances of 3,4-bis(4'-nitrofurazano-3'-yl)furoxan (DNTF) by establishing tannic acid (TA) interface layer on DNTF surface. <i>Chemical Engineering Journal</i> , <b>2022</b> , 434, 134513	14.7	2
11	Crystal structure of 2,4,6,8,10,12-hexanitro-2,4,6,8,10,12-hexaazatetracyclo[5.5.0.0.5.0]dodecane 1/3 hydrate, C <sub>6</sub> H <sub>8</sub> N <sub>12</sub> O <sub>13</sub> . <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , <b>2016</b> , 231, 491-492	0.2	2

10	Design and properties of N,N'-linked bis-1,2,4-triazoles compounds as promising energetic materials. <i>Journal of Molecular Modeling</i> , <b>2020</b> , 26, 130	2	1
9	Molecular dynamics investigation on the morphology of HNIW affected by the growth condition. <i>Journal of Energetic Materials</i> , <b>2019</b> , 37, 44-56	1.6	1
8	Bridged Ti3C2TX MXene Film with Superior Oxidation Resistance and Structural Stability for High-Performance Flexible Supercapacitors. <i>ACS Applied Energy Materials</i> , <b>2022</b> , 5, 2898-2908	6.1	1
7	Investigation of the effect of the CAB/A3 system on HNIW-based PBXs using molecular dynamics. <i>Journal of Molecular Modeling</i> , <b>2018</b> , 24, 186	2	0
6	A novel ternary energetic compound: DAF/DNP/H2O cocrystal. <i>Journal of Energetic Materials</i> , 1-13	1.6	0
5	Reactive molecular dynamics simulation of thermal decomposition for nano-FOX-7. <i>Applied Physics A: Materials Science and Processing</i> , <b>2021</b> , 127, 1	2.6	0
4	Shock Initiation Investigation of a Pressed Trinitrotoluene Explosive. <i>Propellants, Explosives, Pyrotechnics</i> , <b>2021</b> , 46, 1717	1.7	0
3	Crystal structure of hexaaquamagnesium(II) 5,5'-bitetrazole-1,1'-diolate, C <sub>2</sub> H <sub>12</sub> N <sub>8</sub> O <sub>8</sub> Mg. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , <b>2016</b> , 231, 305-306	0.2	
2	Crystal structure of tetraqua((E)-4,4'-(diazene-1,2-diyl)bis(5-oxo-4,5-dihydro-1,2,4-triazol-1-ide)-2N:O)barium(II), C <sub>4</sub> H <sub>10</sub> N <sub>8</sub> O <sub>6</sub> Ba. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , <b>2016</b> , 231, 503-504	0.2	
1	The influence of temperature environmental on performance of HNIW/FOX-7 based PBXs.. <i>Scientific Reports</i> , <b>2022</b> , 12, 4988	4.9	