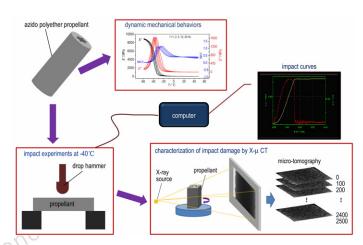
Graphical Abstract

# Impact Damage and Dynamic Mechanical Behaviors of Azido Polyether Propellant at Low Temperature



ZHENG Qi-long, LIU Hai-tao, HU Yi-wen, JIAN Xiao-xia, XIAO Le-qin, ZHOU Wei-liang

Chinese Journal of Energetic Materials, 2017, 25(5): 354-359

The impact load was applied to azido polyether propellant by the instrumented drop hammer device at -40 °C. The impact damage of the sample was characterized by X-ray micro-tomography (X- $\mu$ CT), and the dynamic mechanical properties of the propellant sample were studied by dynamic mechanical analyzer at different loading frequencies.

## Study on the Curing Process of HTPB/IPDI Binder System under Different Catalysts Conditions

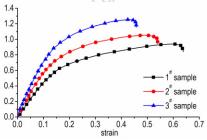
90000 - DBTDL 80000 HTPB-IPDI 70000 TECH 60000 DABCO 50000 TPB nano-ZnO 40000 30000 20000 60 120 180 240 300 360 420 curing time / min

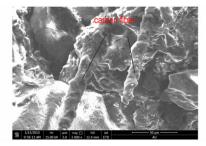
 $\ensuremath{\mathsf{MA}}$  Hui, LIU Yu-cun, CAI Tao, GUO Jia-hu, HU Tuo-ping, LUO Jin, ZHANG Jun

Chinese Journal of Energetic Materials, 2017, 25(5): 360-365

The effects of different catalysts on the curing process of HTPB/IPDI binder system were studied by rheological method. The curing mechanism of HTPB/IPDI system was discussed.

Effect of Short Carbon Fibers on Mechanical Properties of AP/HTPB Base Bleed Propellant 1.41





LIU Zhi-lin, YAO Wen-jin, WANG Xiao-ming, LI Wen-bin, GAO Xiang

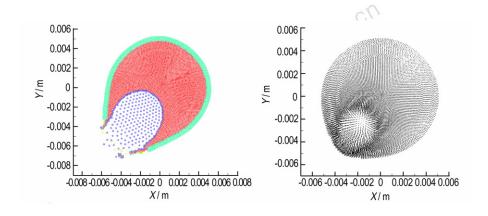
www.energetic-

Chinese Journal of Energetic Materials ,2017 ,25(5): 366-371

The new AP/HTPB base bleed propellants containing different mass fraction of 2 mm short carbon fibers were proposed. The static uniaxial compression tests and tensile tests of new AP/HTPB base bleed propellants were conducted to study the mechanical properties. The effects of short carbon fibers content on the uniaxial compression and tensile properties were studied and the fiber reinforcement mechanism was discussed. The SEM analysis of fracture surface of test samples are also carried out to understand the short carbon fiber distribution in HTPB matrix and its bonding property.

 ${\rm I\hspace{-.1em}I}$ Graphical Abstract

Simulation on Single Gelled Fuel Droplet's Micro Explosion Process Based on SPH Method with Fully Variable **Smoothing Lengths** 

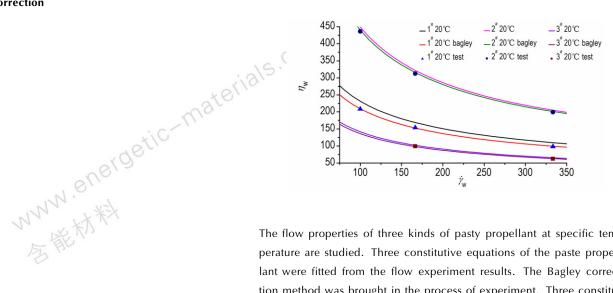


QIANG Hong-fu, ZHANG Lin-tao, CHEN Fu-zhen, LIU Hu, SHI Chao

Chinese Journal of Energetic Materials, 2017, 25(5): 372-378

Numerical simulation study of the micro-explosion process of JP-8 metally gelled fuel droplet was carried out by smoothed SPH method. The changing rule of morphology and physical quantity of the droplet in the micro-explosion process of single gelled fuel droplet was explored. The elastic gelatinizer membrane was simulated by the modified Johnson-Cook Damage (JCD) strength model. The growth, expansion, broken shell, jet etc. processes for internal bubble in the micro-explosion process of gelled droplet were simulated.

### Constitutive Equation of Paste Propellant Based on **Bagley Correction**



LIU Cheng-hao, FENG Feng

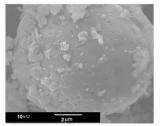
Chinese Journal of Energetic Materials, 2017, 25(5): 379-383

The flow properties of three kinds of pasty propellant at specific temperature are studied. Three constitutive equations of the paste propellant were fitted from the flow experiment results. The Bagley correction method was brought in the process of experiment. Three constitutive equations were refitted according to the Bagley correction parameters, and compared with the previous equations.

Graphical Abstract

## Preparation and Characterization of Metalized Explosive

#### Containing B and Al Powder







SONG Qing-guan, GAO Da-yuan, ZHENG Bao-hui, LI Jing-ming, CAO Wei, CAO Luo-xia, WEN Shang-gang, TAN Kai-yuan

Chinese Journal of Energetic Materials ,2017 ,25(5): 384-390

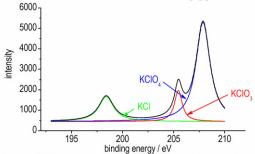
A new HMX based metalized explosive formulation containing B and Al powder, oxidizer AP, polyurethane binder was designed and prepared. The impact sensitivity, friction sensitivity, electrostatic spark sensitivity, cap initiation sensitivity, ignition and propagation properties were measured.

Synthesis and Property Prediction of N-(1,4,6)-trinitrohexahydroimid- azo[4,5-d] imidazole-2 (1H)-ylidene) nitramide

N-(1,4,6-trinitrohexahydroimidazo[4,5-d]imidazol-2(1H)-ylidene) nitramide(TNINA) was synthesized by three stages of nitration reaction using hexahydroimidazo[4,5-d]imidazol-2(1H)-imine as raw material. The optimal condition of the third nitration reaction was obtained and the therrnal stability of target product was studied by thermogravimetry(TG) and differential scanning calorimetry(DSC). The detonation properties were predicted by Kamlet-Jacobs formula.

ZHANG Jun-jun, SHEN Cheng, WANG Peng-cheng, LU Ming *Chinese Journal of Energetic Materials*, 2017, 25(5): 391–395

Effect of Zirconium Powder on Thermal Decomposition of KClO<sub>4</sub> 6000-



2500 2000 KCIO KCIO KCIO 2005 210 binding energy / eV

SUN Ya-lun, LIU Lu, REN Hui, JIAO Qing-jie

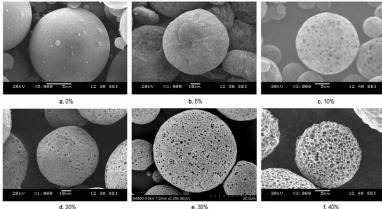
Chinese Journal of Energetic Materials ,2017 ,25(5) : 396-402

The best proportion of Zr and  $\mathrm{KCIO_4}$  were determined through thermodynamic calculation and the influence of Zr on the decomposition process of  $\mathrm{KCIO_4}$  were obtained by XPS. Through  $\mathrm{TG/DTG}$  data of  $\mathrm{KCIO_4}$  and  $\mathrm{Zr/KCIO_4}$  mixture, Kinetic Triplet, critical temperature of thermal explosion and thermodynamic parameters of activation reaction were calculated.

IV Graphical Abstract

## Structure Controlling of Nitrocellulose Base Ball Propel-

lant with Micro-pores

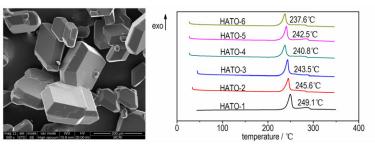


LIN Xiang-yang, SHAO Shan, LI Han, SUN Yi-long, ZHENG Wen-fang, PAN Ren-ming

Chinese Journal of Energetic Materials ,2017 ,25(5): 403-408

•The factors effect on particles diameter, surface porosity, particles sphericity were explored, and the method be used to control particles structure was researched.

Influence of Recrystallization Process on the Thermal Properties and Mechanical Sensitivity of Dihydroxylammonium 5,5'-Bistetrazole-1,1'-diolate(HATO)

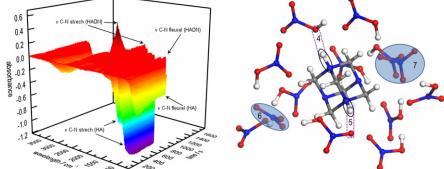


XU Cheng, ZHANG Min, ZHAO Juan, BI Fu-qiang, WANG Ke-yong, ZHU Yan-long, CUI Rong, GE Zhong-xue *Chinese Journal of Energetic Materials*, 2017, 25(5): 409–412

Six kinds of HATO samples with different particle size and morphology were obtained, and characterized by laser granularity instrument and SEM. The thermal decomposition and safety properties of HATO samples were investigated by DSC and impact and friction sensitivity test, respectively.

Salt Formation Mechanism of Urotropine by On-line Infrared Spectroscopy

NNN.8' 各能村

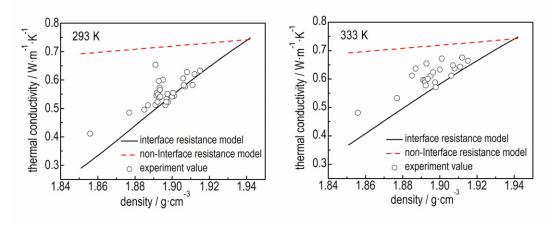


SONG Liang, CHEN Li-zhen, CAO Duan-lin, WANG Jian-long Chinese Journal of Energetic Materials, 2017, 25(5): 413-421

The generation processes of HADN was investigated by on-line infrared spectroscopy. The best on-line infrared spectra were obtained, and the optimization structure of HADN was completed.

V Graphical Abstract

## The Simulation of Thermal Conductivity Coefficien of TATB-based PBX Using the Interface Thermal Resistance Model



ZHOU Xiao-yu, YANG Xue-mei, WEI Xing-wen, WANG Pei Chinese Journal of Energetic Materials, 2017, 25(5): 422-427 The thermal conductivity of TATB-based PBX was predicted by employing the interface-resistance-based thermalcondctivity model. The results of simulation were well agreed with the test values using the flash diffusivity method.

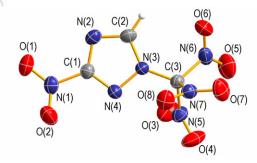
## Research Progress in Energy Conversion Components for **MEMS Initiating Explosive Device**

ZHANG Bin, CHU En-yi, REI Wei, WANG Ke-xuan, LI Hui,

Chinese Journal of Energetic Materials, 2017, 25(5): 428-436

The progressin the energy conversion components for MEMS initiating explosive device in recent years was reviewed with 38 references.

www.energetic-materials. Crystal Structure and Property Prediction of 1-Trinitromethyl-3-nitro-1,2,4-triazole



YIN Xin, MA Qing, WANG Jun, WANG Shu-min

Chinese Journal of Energetic Materials ,2017 ,25(5): 437-440

An energetic material 1-trinitromethyl-3-nitro-1,2,4-triazole (TNMNT) was synthesized via a new synthetic method. Its crystal structure was characterized by X-ray single crystal diffraction for the first time, and the crystal of TNMNT belongs to monoclinic system and space group is  $P2_1/c$ .

Executive editor: GAO Yi ZHANG Qi WANG Yan-xiu JIANG Mei