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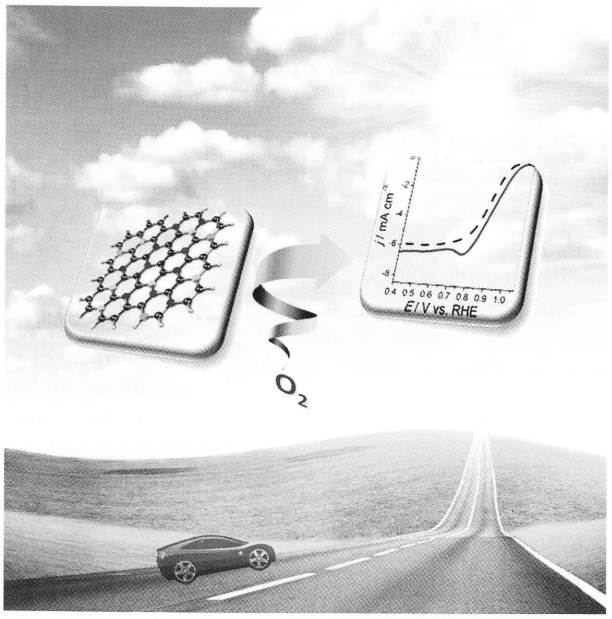
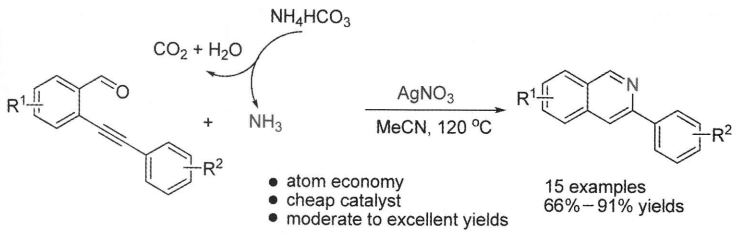
## COVER PICTURE

**The cover picture shows** multiple elements doped reduced graphene oxide (rGO) as an efficient electrocatalyst for the oxygen reduction reaction (ORR). As alternatives to Pt-based electrocatalysts, the development of nonprecious metal catalysts with high performance in the cathodic ORR is highly desirable for large-scale commercialization of automotive fuel cells. A simple two-step doping method for the preparation of pentabasic (Fe, B, N, S, P)-doped rGO as an efficient ORR catalyst has been developed. Simultaneously doping appropriate amounts of Fe and B into the ternary doped N, S, P-rGO composite produced a synergistic effect for enhancing ORR activity, which endowed the prepared catalyst with even higher ORR performance than that of the commercial Pt/C catalyst in alkaline electrolytes. More details are discussed by Liu *et al.* on page 878—886.




## COMMUNICATIONS






An operationally simple approach for the tandem synthesis of isoquinolines by the reaction of *o*-alkynylaldehydes with ammonium bicarbonate *via* Ag-catalyzed 6-endo-dig ring closure is described. The reaction conditions and the scope of the reaction are examined, and a variety of substituted isoquinolines are prepared in moderate to excellent yields.

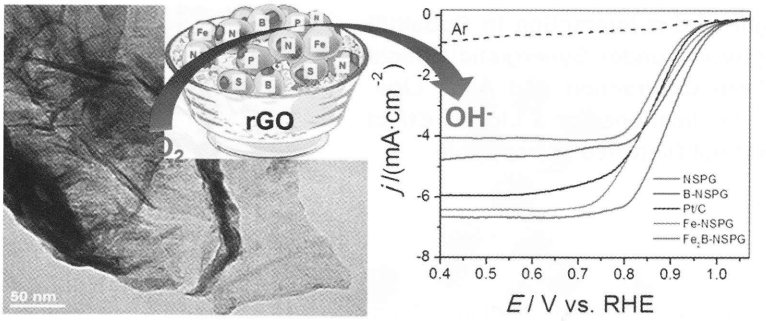
# CONTENT



## *FULL PAPERS*

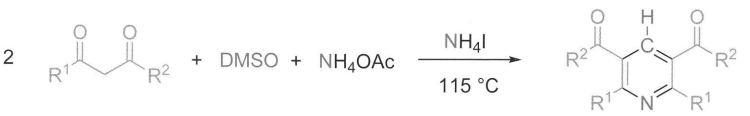




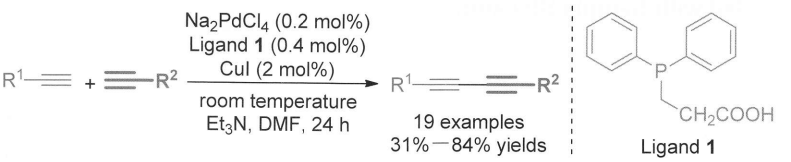

A simple two-step doping method for the preparation of pentabasic (Fe, B, N, S, P)-doped rGO as an efficient ORR catalyst has been developed. Simultaneously doping appropriate amounts of Fe and B into the ternary doped N, S, P-rGO composite produced a synergistic effect for enhancing ORR activity, which endowed the prepared catalyst with even higher ORR performance than that of the commercial Pt/C catalyst in alkaline electrolytes.



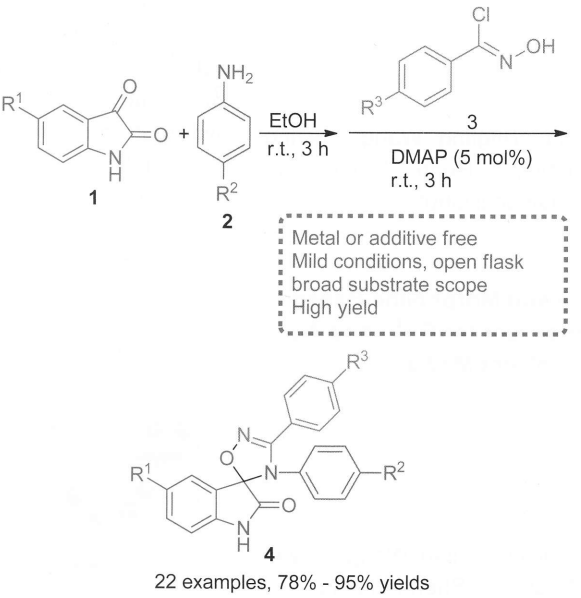



A one-pot synthesis of Hantzsch pyridines was achieved through $NH_4I$-promoted condensation of 1,3-dicarbonyl compounds with DMSO and $NH_4OAc$, in which the C4 of the pyridine rings was derived from the decomposition of DMSO.







# CONTENT



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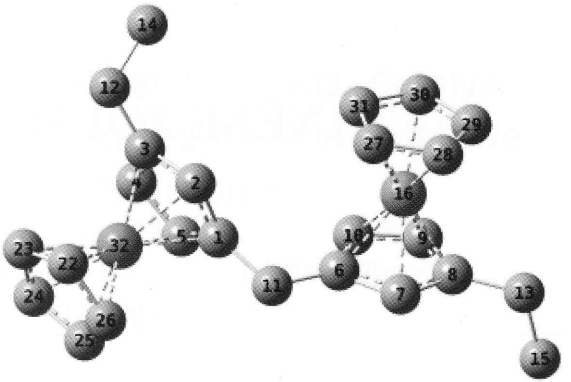
**Correlation between Molecular Structure of Ferrocene Derivative and Impact Sensitivity of Fine-AP/Ferrocene Derivative Mixture**




This figure was the equilibrium molecular configuration of DEFM, one of the new effective DFDs. The bond lengths, the electrostatic potentials of characteristic groups and the DFD molecules, and the molecular frontier orbital energy of DFDs were calculated. The results showed that the DFDs with methylene bridge group were more stable than those with methylethylidene bridge group. The impact sensitivities of the fine-AP/DFD mixtures were in inverse proportion to the stabilities of the DFDs.

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