

## Notes to the Editor

doublet corresponding to a free amine at 3315 and 3395  $\text{cm}^{-1}$  and the absorption corresponding to the nitro group disappeared. The n.m.r. spectra of the diamine (III) showed a singlet at 1.6 $\delta$  corresponding to 6 aliphatic hydrogens, a singlet at 2.75 $\delta$  corresponding to 4 amine hydrogens and a multiplet at 6.8–7.2 $\delta$  corresponding to 6 aromatic hydrogens. From the mass spectrum of the diamine (III), the molecular weight was also checked and found to be 410.

The diamine (III) was dissolved in  $\text{CHCl}_3$ , sodium carbonate solution and sodium lauryl sulphate solution added and the content of the flask stirred vigorously for 5 min. All of the terephthaloyl chloride in chloroform was added and stirring was continued for another 5 min. Hexane was then added and the contents were slowly stirred for 2 min, filtered, washed with water, chloroform and hexane. The residue was redissolved in DMF and reprecipitated with methanol and dried at 100°C under reduced pressure. The i.r. spectrum of the polymer (IV) showed the disappearance of the free amine frequency and the appearance of peaks at 1595 and 1645  $\text{cm}^{-1}$  cor-

responding to amide  $-\text{NH}$  stretching and amide  $\text{C}=\text{O}$  stretching respectively and a broad band at 3200  $\text{cm}^{-1}$  again corresponding to  $\text{N}-\text{H}$  stretching, confirming the structure (IV) for the polymer<sup>12</sup>. The polymer was soluble in dimethyl sulphoxide and dimethylformamide. The inherent viscosity was found to be 1.7 in DMF at 30°C at a concentration of 0.025 g/100 ml. Differential thermal analysis showed that the polymer was unaffected up to 380°C. Further work regarding molecular weight, viscosity etc. are in progress.

## ACKNOWLEDGEMENTS

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## REFERENCES

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## ERRATA

'Some aspects of stereoregulation in the stereospecific polymerization of vinyl monomers' by Piero Pino and Ulrich W. Suter, *Polymer* 1976, 17, 977–995

Page 978, left hand column, line 28: *for* 33, *read* 33 monomeric units  
Page 979, Table 1, 2 lines from bottom, under Solvent: *for* pentene, *read* pentane  
Page 981, Table 2, line 4, under Dyads (m): *for* 0.8 *read* 0.3; line 19, under Monomers: *for* Menthyl vinyl ketone *read* methyl vinyl ketone  
Page 982, left hand column line 4: *for* include *read* induce; right hand column line 23: *for* anylsodium *read* amylsodium  
Page 983, Table 4 2 lines from bottom under Stereoregularity: *for* Atactic, (m) = 40, *read* Atactic, (m) = 0.40; bottom line, under Stereoregularity: *for* predicted, *read* predominantly  
Page 984, right hand column, line 34 should read: With ethyl-(Z)- $\beta$ - $d_1$ -methacrylate, Fowells, Schuerch, Bovey and Hood . . .  
Page 985, Table 9, footnote: *for* (E)- $\beta$ - $d_1$ -methyl methacrylate, *read* methyl-(Z)- $\beta$ - $d_1$ -methacrylate  
Page 986, Table 10, caption and left

hand column, line 27: *for* (Z)- $\alpha$ -ethyl- $\beta$ - $d_1$ -methyl acrylate, *read* ethyl-(Z)- $\beta$ - $d_1$ -methacrylate; left hand column, line 32: *for* contact with ion pairs, *read* contact ion pairs; left hand column, line 40 should read: For poly(vinyl ethers), for instance, a ring structure is suggested as shown in Figure 3, the oxygenated chain acting as a specific solvating agent itself<sup>99</sup>; right hand column, line 6: *for* the same *read* some  
Page 987, Table 11, bottom line under Monomer: *for* 1-Methyl propyl vinyl, *read* 1-methyl propyl vinyl ether; left hand column line 3: *for* (Z)- $\alpha$ -ethyl- $\beta$ - $d_1$ -methacrylate, *read* ethyl-(Z)- $\beta$ - $d_1$ -methacrylate  
Page 988, schemes (3) and (4): *for* Me, *read* [Me]; right hand column, 5 lines from the bottom: *for* (2b), *read* (4b)  
Page 989, left hand column, 10 lines from the bottom: *for* (E)-1-deuteropropylene, *read* (E)-1,3,3,3- $d_4$ -propylene  
Page 990, left hand column, line 7: *for*

(equation 2), *read* (equation 11); right hand column, scheme (9): *for*  $\text{Ti}(\text{Bz})_4/\text{Al}(\text{Br})_3$ , *read*  $\text{Ti}(\text{Bz})_4/\text{Al}(\text{Bz})_3$   
Page 992, Figure 6 caption: *for*  $H_M - H_R = 2.2$  kcal/mol;  $S_M - S_R = 5.0$  cal/mol K, *read*  $H_i - H_s = 2.2$  kcal/mol;  $S_i - S_s = 5.0$  cal/mol K; Scheme 12, structure (1), the geminal 'CH<sub>3</sub>' and 'H' on the propylene complexed to the catalyst should be interchanged  
Page 993, reference 19 should read: White, D. M. *J. Am. Chem. Soc.* 1960, 82, 5678; Brown, Jr J. F. and White, D. M. *J. Am. Chem. Soc.* 1960, 82, 5671  
Page 994 reference 46: *for* Kirsch Yu. E., *read* Kirsh, Yu. E.; reference 68: *for* Isio, G. A., *read* Audisio, G.; references 61, 77 and 86 *for* Urgu, T. *read* Uryu, T.  
Throughout the text *for*  $4(\text{mm})(\text{rr})(\text{mr})^2$  *read*  $(\text{mm})(\text{rr})(\text{mr})^2$   
We apologise for the omission of these printers and authors errors at the proof stage.