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Re-entrance of Poly(N,N-diethylacrylamide) in $\mathrm{D}_{2} \mathrm{O} /$ d-Ethanol Mixture at $27{ }^{\circ} \mathbf{C} .{ }^{1}$ HE CHENG, Institute of High Energy Physics CAS - The reentrance of poly( $\mathrm{N}, \mathrm{N}$-diethylacrylamide) (PDEA) in $\mathrm{D}_{2} \mathrm{O} /$ d-ethanol mixtures (i.e., the coil-to-spherical aggregates-to-coil transition) has been observed at $27^{\circ} \mathrm{C}$ by small-angle neutron scattering (SANS). PDEA has a lower critical solution temperature (LCST) phase diagram in the $\mathrm{D}_{2} \mathrm{O}$ rich region and is soluble in the $\mathrm{D}_{2} \mathrm{O}$-poor region for all of the observed temperature ranges. Its spinodal temperature decreases first from $33.5^{\circ} \mathrm{C}$ in pure $\mathrm{D}_{2} \mathrm{O}$ to $26.7^{\circ} \mathrm{C}$ in $80 \% \mathrm{D}_{2} \mathrm{O} / 20 \%$ d-ethanol and then increases to $283.1^{\circ} \mathrm{C}$ in $50 \% \mathrm{D}_{2} \mathrm{O} / 50 \%$ d-ethanol With the further decrease of $\mathrm{D}_{2} \mathrm{O}$ content, PDEA dissolves well, and its phase boundary can no longer be observed by SANS. The ternary random phase approximation model (RPA) is used to analyze the SANS profiles, and three Flory-Huggins interaction parameters ( $\chi_{\text {PDEAdethanol }}$, $\chi_{\text {PDEAD2O }}$ and $\chi_{\text {dethanolD2O }}$ ) are obtained. When a small amount of d-ethanol is added to the system, it has a strong interaction with $\mathrm{D}_{2} \mathrm{O}$, so it directly gets distributed into the water structure and makes a negative contribution to the dissolution of PDEA ( $\chi_{\text {dethanolD2O }}$ is much smaller than $\chi_{\text {PDEAdethanol }}$ and $\chi_{\text {PDEAD2O }}$ ). Neither d-ethanol nor $\mathrm{D}_{2} \mathrm{O}$ wants to help the dissolution of PDEA in the first place, until the structure of mixed solvents tends to be pure d-ethanol in the $\mathrm{D}_{2} \mathrm{O}$-poor region.
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