実験報告書様式(一般利用課題·成果公開利用)

(※本報告書は英語で記述してください。ただし、産業利用課題として採択されている方は日本語で記述していただいても結構です。)

Experimental Report	承認日 Date of Approval 承認者 Approver 提出日 Date of Report
課題番号 Project No.	装置責任者 Name of Instrument scientist
2017A0002	Jun-ichi Suzuki
実験課題名 Title of experiment	装置名 Name of Instrument/(BL No.)
Does Aggregation of Ethanol Molecules in Imidazolium-based	TAIKAN/(BL15)
Ionic Liquids Affect the Complex Formation of Metal Ions?	実施日 Date of Experiment
実験責任者名 Name of principal investigator	2017 Nov. 11 10:00–Nov. 14 10:00
Toshiyuki Takamuku	
所属 Affiliation	
Saga University	

試料、実験方法、利用の結果得られた主なデータ、考察、結論等を、記述して下さい。(適宜、図表添付のこと)

Please report your samples, experimental method and results, discussion and conclusions. Please add figures and tables for better explanation.

1. 試料 Name of sample(s) and chemical formula, or compositions including physical form.

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In the present machine time for 72 h, SANS experiments were conducted on the following systems.
System A:
Imidazolium-based ionic liquids-ethanol (EtOH) solutions at 25 °C
1. C<sub>6</sub>mimTFSA-Ethanol-d<sub>6</sub> x<sub>FtOHd6</sub> = 0.85, 0.90, 0.925, 0.95, 0.97, 0.99
System B:
Imidazolium-based ionic liquids-dimethyl sulfoxide (DMSO) solutions at 25 °C
1. C_n mimTFSA-DMSO-d_6 with n = 2-12 at x_{DMSOd_6} = 0.95, 0.97
System C:
Imidazolium-based ionic liquids-acetonitrile (AN) solutions at 25 °C
1. C_2mimTFSA-Acetonitrile-d_3 = 0.80, 0.90, 0.95, 0.99
2. C<sub>8</sub>mimTFSA-Acetonitrile-d<sub>3</sub> x<sub>ANd3</sub> = 0.80, 0.90, 0.95, 0.99
System D:
Temperature dependence of the inhomogeneity of Imidazolium-based ionic liquids-1,4-dioxane (DIO) solutions
1. C<sub>8</sub>mimTFSA-1,4-Dioxane-d_8 x_{DIOd8} = 0.9825 in the temperature range of 50-31 °C
System E:
Lysozyme-alcohol-water solutions at 25 °C
1. Lys/2-PrOH-d_8-D<sub>2</sub>O-HFIP x_{2-PrOHaB} = 0.3, 0.35
2. Lys/2-PrOH-d_8-D<sub>2</sub>O/H<sub>2</sub>O-HFIP x_{2-PrOH} = 0.3, 0.35
3. Lys/D<sub>2</sub>O-HFIP x_{\text{HFIP}}= 0.2
System F:
1. D_2O-3-methylpyridine (fixed at 9 vol.%) + NaBPh<sub>4</sub> (0 to 85 mM) + PPh<sub>4</sub>Cl (0 to 40 mM)
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2. 実験方法及び結果(実験がうまくいかなかった場合、その理由を記述してください。)

Experimental method and results. If you failed to conduct experiment as planned, please describe reasons.

Systems A-C:

The SANS measurements for Systems A-C were made to clarify the relation between the inhomogeneity of the ionic liquids (ILs)-molecular liquids (MLs) solutions and the complex formation of Ni²⁺ in their binary solutions. The MLs includes EtOH, DMSO, and AN.

From the present SANS experiments, together with the previous ones (2015A0038, 2016A0008, 2017A0006), the mixing states of the C_nmimTFSA-ML solutions with n = 2-12 can be clarified at the mesoscopic level. The DMSO and AN solutions did not show significant SANS intensities. Thus, both MLs are homogeneously mixed with the imidazolium-based ILs. On the contrary, the SANS intensities of the EtOH solutions were recorded on the TAIKAN spectrometer. The Ornstein-Zernike correlation lengths ξ were estimated for all of the C_nmimTFSA-Ethanol-*d*₆ solutions and plotted as a function of ethanol mole fraction x_{EtOHd6} in Fig. 1. As seen, the maximum of the ξ appears at the higher x_{EtOHd6} with elongating the IL alkyl chain n. Moreover, the maximum ξ decreases with the alkyl chain length. The dispersion force between the EtOH ethyl group and the IL alkyl group may influence the mixing states of the IL-ML solutions. On the basis of the SANS results, we will discuss the effects of the mixing states of the IL-ML solutions on the complex formation of Ni²⁺ in their solutions.



Fig. 1 Ornstein–Zernike correlation lengths ξ of C_n mimTFSA–Ethanol– d_6 solutions as a function of ethanol mole fraction.

System D:

We found that C_8 mimTFSA-DIO solutions separate into the IL-rich and DIO-rich phases with lowering temperature. As shown in the phase diagram (Fig. 2), the solutions show the upper critical solution temperature (UCST) at 31.6 °C and $x_{DIO} = 0.9825$. In the present experiments, we observed SANS profiles of the C_8 mimTFSA-DIO solution at $x_{DIO} = 0.9825$ with lowering temperature toward the UCST. Fig. 3 shows the Ornstein-Zernike correlation lengths ξ estimated for the solution as a function of temperature. The ξ remarkably increases with lowering temperature. This suggests that the inhomogeneity of the solution enhances as the temperature decreases towards the UCST. Moreover, the plots of the ξ values against the reciprocal temperature showed that phase separation of the C_8 mimTFSA-DIO solution takes place in the 3D-Ising manner. This is the first result on the mechanism of phase separation of IL-ML solutions.





Fig. 2 Phase diagram of C_n mimTFSA-DIO solutions with n = 2, 4, and 8.



Fig. 3 Ornstein-Zernike correlation lengths ξ of C₈mimTFSA-DIO solution at x_{DIO} = 0.9825.

System E:

In the present and previous machine times (2015A0038, 2016A0008, 2017A0006), we made the SANS measurements at 25 °C on lysozyme-alcohol-water solutions using the TAIKAN spectrometer to elucidate the change in the 3D structure of lysozyme depending on the alcohols and their mole fraction. Here, the alcohols are 2-propanol (2-PrOH) and 1,1,1,3,3,3-hexafluoroisopropanol (HFIP). The SANS profiles of the 2-PrOH solutions did not significantly change with increasing x_{2-PrOH} from 0 to 0.3. Thus, the 3D structure of lysozyme is not affected by 2-PrOH in the mole fraction range. In contrast, the SANS intensities of the HFIP solutions strengthened with increasing x_{HFIP} from 0 to 0.05. With further increasing x_{HFIP} to 0.2, the intensities decrease. Therefore, the 3D structure of lysozyme obviously changes with increasing HFIP content. This finding agrees with the results of 2D structure by circular dichroism (CD) spectroscopy. We are analyzing the SANS profiles on the 3D structure of lysozyme in aqueous solutions will be clarified on the basis of the SANS profiles of HFIP on the 3D structure of lysozyme in aqueous solutions will be clarified on the basis of the SANS profiles using the set of the CD and IR spectroscopy.

System F:

In the present and previous machine times (2017A0006), the feature of the lamellar structure formed in System F was investigated by changing the concentration of NaBPh₄ and PPh₄Cl. One example of the SANS results is shown in Fig. 4. Here, the concentration of NaBPh4 was fixed at 85 mM. With increasing the amount of PPh₄Cl, the Bragg peak around Q = 0.03 Å⁻¹ disappears, while the intensity at low-Q region increases. In this manner, it was shown that the addition of PPh₄Cl decreases the regularity of the membrane, however, induces the micrometer-sized ordered structures. This finding is an important clue to clarify how the charged substance affects the membrane composed of organic solvent and antagonistic salts.



