



Volumetric Studies of Sodium Dodecyl Sulphate in Aqueous and Aqueous Amino Acid Solutions

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ABSTRACT

The volumetric studies on the binary aqueous solutions of sodium dodecyl sulphate (SDS) and ternary SDS-amino acid-water systems have been carried out as a function of concentration and at 15, 20, 25, 30, 35, 40, 45 and 50 °C. In this study the variation of apparent molar volume (ϕ_v) of SDS was observed in the low concentration region. The limiting apparent molar volume (ϕ_v^0) values of SDS in aqueous amino acid solutions are greater than those in water and increase with the concentration at all the studied temperatures. The greater ϕ_v^0 value in amino acid solution has been demonstrated strong ion-ion interactions in the ternary system. Apparent molar expansions (Φ_E) in aqueous solution and ternary SDS-amino acid-water systems are very sensitive to hydrophobic and hydrophilic solute-solute interactions. Expansibility, α and apparent molar expansion, Φ_E reflect from the same structural alteration in the solutions. In our investigation, critical micelle concentration (c.m.c) of SDS at various SDS-amino acid-water systems also determined by apparent molar volume (ϕ_v). Among various physical parameters, apparent molar volume, limiting apparent molar volume, Apparent molar expansion has been recognized as a quantity that is sensitive to structural changes occurring in solutions.

Keywords : Sodium dodecyl sulphate, glycine, alanine, apparent molar volume, apparent molar expansion, critical micelle concentration (c.m.c).

INTRODUCTION

Volumetric studies of SDS in aqueous and in aqueous amino acid provides great help to determine the nature of interactions and properties of this solution. And it is very important because of great applications of SDS in Pharmaceuticals and Biotechnological processes. SDS represents a potentially effective topical microbicide, which can also inhibit and possibly prevent infection by

various enveloped and non-enveloped viruses such as the Herpes simplex viruses, HIV, and the Semliki Forest virus.

In this paper, We reports the changes in the physical properties due to the mixing of water, sodium dodecyl sulphate & amino acid are thought of arising from hydrophobic interactions among solute and solvent molecules. Here has been used ternary process and the results of volumetric studies

on the binary aqueous solutions of sodium dodecyl sulphate(SDS) and ternary SDS-amino acid-water systems gives vast information about volumetric properties. Here we used two amino acids (glycine & dl-alanine) and sodium dodecyl sulphate(SDS). Surfactant molecules (e.g. CTAB, SDS, TritonX-100, etc.) self-aggregate into super molecular structures when dissolved in water or oil. The simplest aggregate of these surfactant molecules is called a micelle; and the dispersion of the aggregates in water or oil is referred to as a micellar solution. A typical micelle has size of $\sim 50 \text{ \AA}$ and is made of about 100 surfactant molecules. The surfactant molecule consists of two parts, namely, a polar hydrophilic head group and an apolar hydrophobic tail (hydrocarbon chain). The concentration above which micelle formation becomes appreciable is termed the *critical micelle concentration* (c.m.c).

Micelle formation is as typical hydrophobic process in water. In aqueous medium, surfactant molecules with their long hydrophobic tails undergo hydrophobic hydration. As the surfactant concentration increases, the association of surfactant molecules occurs by hydrophobic interaction and this result in the removal of the non-polar portion of the molecule from the external aqueous environment to form the interior of the micelle while the hydrophilic groups are exposed to the aqueous environment

Therefore, the present studies are to determined apparent molar volume, apparent molar expansion and critical micelle concentration (c.m.c) at low region concentration in binary and ternary system. A review of literature shows that only a few authors have attempted to determine of these properties in ternary system at low region concentration.

EXPERIMENTAL

The Surfactant used in this study was sodium dodecyl sulphate(SDS). The c.m.c value of SDS in water reported in the literature value closely around 0.009m at 25°C. The values of mean aggregation number of SDS vary from 58 to 64. Glycine (purity. Mass fraction ≥ 0.99), DL-alanine (purity. Mass fraction ≥ 0.99) procured from Fluka Chemical Company, Switzerland were used without

further purification. Supplied distilled water was redistilled and deionized by passing through two ion exchange columns. The deionized water was distilled again in alkaline KMnO_4 medium and used for preparation of solution. Conductivity of this deionized water was found to be $5 \times 10^{-6} \text{ S.cm}^{-1}$. An electric balance with an accuracy of $\pm 0.0001 \text{ g}$ was used for weighting. The densitometer (DSA-5000, Anton Paar, Austria) was used for the measurements of density and ultrasonic sound velocity. The solutions were prepared by weight immediately before the measurement. The ternary solutions were prepared by mixing appropriate mass of the components. The amount of each component was later converted into their mole fraction. Precautions were taken to prevent the introduction of moisture into the experimental examples. Each time, the solution was prepared immediately before the density measurement.

RESULTS AND DISCUSSION

Apparent molar volume

The apparent molar volume, ϕ_v , was calculated from the measured density of the solution using the relation

$$\phi_v / \text{cm}^3 \cdot \text{mol}^{-1} = \frac{1000}{m\rho\rho_0}(\rho_0 - \rho) + \frac{M_2}{\rho} \quad \dots(1)$$

Where M_2 , m , ρ and ρ_0 respectively the molar mass of solute, molality of solution, density of solution and density of solvent. Density data of solution and solvent are given in Table 1 and the calculated values at different temperatures are given in Table 2 as a function of molality of the solutions. The plots of the against are presented in Figs. 1-3.

For surfactant solutions physical properties such as, osmotic pressure, turbidity, electrical conductance and surface tension were used to determine the critical micelle concentration (c.m.c). In our investigation we tried to observe the variation of apparent molar volume of SDS as we focused our study in the low concentration region. The $\phi_v - \sqrt{m}$ plots show sudden change in ϕ_v value at a particular molality after which show linear change. The c.m.c of SDS in aqueous and aqueous amino acid solutions were determined by extrapolation method. The c.m.c values of SDS in aqueous and aqueous

Table.1: Density (ρ) and (ρ_0) of SDS in water and aqueous amino acid solutions as a function of molality of SDS at different temperature.

WATER+SDS		Water, (Density, ρ_0 /g.cm ⁻³)						
m (mol.kg ⁻¹)	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.00000	0.999101	0.998206	0.997045	0.995646	0.994032	0.992215	0.990212	0.988035
	Water +SDS, (Density, ρ_0 /g.cm ⁻³)							
0.000340	0.999138	0.998239	0.997080	0.995679	0.994063	0.992246	0.990243	0.988065
0.000890	0.999171	0.998272	0.997109	0.995708	0.994094	0.992273	0.990271	0.988091
0.001360	0.999189	0.998289	0.997124	0.995724	0.994108	0.992288	0.990284	0.988106
0.014990	0.999932	0.999007	0.997823	0.996400	0.994766	0.992932	0.990917	0.988723
0.029990	1.000638	0.999694	0.998491	0.997054	0.995406	0.993558	0.991528	0.989326
0.043980	1.001294	1.000330	0.999111	0.997660	0.995999	0.994138	0.992095	0.989885
0.058070	1.001942	1.000959	0.999725	0.998258	0.996580	0.994706	0.992655	0.990433
0.071994	1.002582	1.001580	1.000327	0.998845	0.997155	0.995273	0.993208	0.990979
0.085940	1.003218	1.002201	1.000932	0.999436	0.997731	0.995835	0.993759	0.991518
0.100850	1.003897	1.002860	1.001571	1.000062	0.998345	0.996433	0.994348	0.992095
0.114980	1.004497	1.003443	1.002140	1.000615	0.998886	0.996965	0.994868	0.992606
0.128601	1.005091	1.004022	1.002706	1.001166	0.999424	0.997491	0.995384	0.993111
0.142570	1.005672	1.004584	1.003252	1.001702	0.999945	0.998002	0.995885	0.993605
0.157000	1.006295	1.005189	1.003841	1.002276	1.000507	0.998549	0.996421	0.994132
0.171940	1.007184	1.005953	1.004580	1.002988	1.001199	0.999222	0.997076	0.994769
0.186100	1.007654	1.006463	1.005084	1.003485	1.001687	0.999707	0.997556	0.995244
0.198450	1.008073	1.006918	1.005524	1.003916	1.002109	1.000121	0.997964	0.995647

WATER+GLYCINE(m=1.3)+SDS

m (mol.kg⁻¹)	Water+Glycine, (Density, ρ_d/g.cm⁻³)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
1.300000	1.040102	1.038734	1.037184	1.035465	1.033588	1.031560	1.029387	1.027090		
	Water+Glycine+SDS, (Density,ρ/g.cm⁻³)									
0.001592	1.040001	1.038632	1.037080	1.035363	1.033484	1.031455	1.029286	1.026989		
0.014431	1.040170	1.038785	1.037225	1.035493	1.033600	1.031558	1.029370	1.027055		
0.013725	1.040582	1.039199	1.037631	1.035897	1.034005	1.031961	1.029793	1.027481		
0.027987	1.041098	1.039701	1.038116	1.036369	1.034462	1.032411	1.030222	1.027912		
0.039750	1.041758	1.040340	1.038746	1.036985	1.035071	1.033010	1.030810	1.028500		
0.053373	1.042167	1.040736	1.039128	1.037357	1.035429	1.033359	1.031147	1.028815		
0.069147	1.042634	1.041187	1.039562	1.037775	1.035839	1.033752	1.031531	1.029188		
0.085642	1.043342	1.041874	1.040231	1.038427	1.036474	1.034375	1.032142	1.029785		
0.100063	1.043642	1.042165	1.040509	1.038694	1.036731	1.034623	1.032380	1.030013		
0.114261	1.044118	1.042627	1.040959	1.039131	1.037151	1.035034	1.032781	1.030405		
0.127985	1.044650	1.043143	1.041462	1.039622	1.037633	1.035503	1.033238	1.030860		
0.140583	1.045041	1.043519	1.041828	1.039977	1.037976	1.035838	1.033566	1.031176		
0.156884	1.045840	1.044300	1.042584	1.040720	1.038705	1.036551	1.034267	1.031865		
0.171385	1.046048	1.044496	1.042772	1.040892	1.038869	1.036707	1.034413	1.032001		
0.185656	1.046715	1.045145	1.043406	1.041512	1.039476	1.037304	1.035001	1.032578		
0.195929	1.046825	1.045245	1.043497	1.041600	1.039554	1.037372	1.035066	1.032637		

WATER+GLYCINE(0.883)+SDS

m (mol.kg ⁻¹)	Water+Glycine, (Density, ρ _g .cm ⁻³)							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.883000	1.027347	1.026115	1.024672	1.023047	1.021241	1.019274	1.017154	1.014884
	Water+Glycine+SDS, (Density,ρ_g.cm⁻³)							
0.000680	1.026967	1.025736	1.024298	1.022680	1.020879	1.018903	1.016781	1.014521
0.001544	1.026991	1.025757	1.024318	1.022690	1.020888	1.018921	1.016801	1.014534
0.014193	1.027606	1.026358	1.024903	1.023261	1.021448	1.019469	1.017330	1.015050
0.024522	1.028213	1.026949	1.025483	1.023827	1.022001	1.020011	1.017861	1.015592
0.037557	1.028653	1.027376	1.025895	1.024227	1.022389	1.020389	1.018239	1.015949
0.051069	1.029154	1.027860	1.026365	1.024685	1.022835	1.020826	1.018665	1.016366
0.071936	1.030103	1.028781	1.027256	1.025562	1.023689	1.021660	1.019490	1.017178
0.086225	1.030621	1.029288	1.027754	1.026040	1.024159	1.022120	1.019932	1.017609
0.100026	1.031096	1.029747	1.028199	1.026474	1.024579	1.022529	1.020331	1.017998
0.114767	1.031727	1.030363	1.028797	1.027057	1.025150	1.023087	1.020881	1.018536
0.128215	1.032291	1.030910	1.029330	1.027575	1.025657	1.023583	1.021365	1.019012
0.142658	1.032819	1.031421	1.029829	1.028061	1.026131	1.024046	1.021819	1.019455
0.157116	1.033287	1.031872	1.030267	1.028486	1.026543	1.024450	1.022210	1.019839
0.171290	1.033799	1.032367	1.030754	1.028962	1.027006	1.024799	1.022550	1.020171
0.185205	1.034290	1.032852	1.031215	1.029392	1.027429	1.025310	1.022990	1.020599
0.198478	1.034735	1.033277	1.031632	1.029814	1.027838	1.025711	1.023443	1.021046

WATER+DL-ALANINE(1.216)+SDS

m (mol.kg ⁻¹)	Water+DL-alanine (Density, ρ , g.cm ⁻³)							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
1.216000	1.033900	1.032648	1.031192	1.029555	1.027746	1.025772	1.023677	1.021416
Water+DL-alanine+SDS, (Density, ρ, g.cm⁻³)								
0.001430	1.033233	1.031988	1.030539	1.028904	1.027097	1.025127	1.023006	1.020745
0.013983	1.034105	1.032836	1.031368	1.029717	1.027895	1.025912	1.023779	1.021511
0.026784	1.034966	1.033681	1.032196	1.030528	1.028692	1.026695	1.024553	1.02227
0.041951	1.035471	1.034171	1.032670	1.030988	1.029137	1.027130	1.024962	1.022663
0.055598	1.035995	1.034676	1.033160	1.031465	1.029601	1.027582	1.025414	1.023115
0.070241	1.036521	1.035187	1.033654	1.031943	1.030068	1.028036	1.025859	1.023547
0.082870	1.037049	1.035699	1.034151	1.032431	1.030542	1.028501	1.026312	1.023993
0.098097	1.037568	1.036201	1.034639	1.032900	1.031001	1.028949	1.026749	1.024419
0.108665	1.037649	1.036272	1.034703	1.032958	1.031052	1.028989	1.026784	1.024460
0.125733	1.038211	1.036816	1.035228	1.033468	1.031548	1.029474	1.027257	1.024906
0.140410	1.038999	1.037585	1.035980	1.034206	1.032267	1.030181	1.027948	1.025589
0.153293	1.039494	1.038064	1.036445	1.034656	1.032708	1.030609	1.028371	1.026003
0.168333	1.039877	1.038437	1.036807	1.035008	1.033052	1.030946	1.028695	1.026318
0.182099	1.040194	1.038737	1.037096	1.035282	1.033317	1.031200	1.028944	1.026557
0.198069	1.040507	1.039040	1.037383	1.035559	1.033581	1.031456	1.029186	1.026790

WATER+DL-ALANINE(0.827)+SDS

m (mol.kg ⁻¹)	Water+DL-alanine (Density, ρ ₀ /g.cm ⁻³)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
0.827000	1.022989	1.021845	1.020478	1.018912	1.017159	1.015229	1.013160	1.010920		
	Water+DL-alanine+SDS, (Density,ρ/g.cm⁻³)									
0.000760	1.023066	1.021921	1.020553	1.018986	1.017233	1.015303	1.013230	1.010990		
0.001556	1.023067	1.021921	1.020552	1.018984	1.017230	1.015298	1.013226	1.010985		
0.014299	1.023563	1.022400	1.021014	1.019432	1.017664	1.015725	1.013631	1.011383		
0.028638	1.024186	1.023003	1.021602	1.020002	1.018222	1.016272	1.014165	1.011906		
0.042769	1.024702	1.023502	1.022086	1.020474	1.018681	1.016717	1.014600	1.012333		
0.055985	1.025339	1.024121	1.022690	1.021064	1.019259	1.017283	1.015157	1.012875		
0.070499	1.025950	1.024723	1.023279	1.021635	1.019816	1.017831	1.015690	1.013399		
0.083252	1.026449	1.025200	1.023736	1.022084	1.020254	1.018258	1.016107	1.013811		
0.099087	1.027024	1.025757	1.024276	1.022609	1.020763	1.018753	1.016593	1.014286		
0.112982	1.027574	1.026293	1.024799	1.023115	1.021259	1.019239	1.017066	1.014749		
0.127136	1.027840	1.026543	1.025036	1.023342	1.021475	1.019444	1.017263	1.014939		
0.139751	1.028550	1.027234	1.025713	1.024003	1.022122	1.020080	1.017889	1.015556		
0.154131	1.028734	1.027408	1.025876	1.024141	1.022235	1.020217	1.018018	1.015665		
0.169950	1.029680	1.028332	1.026778	1.025043	1.023137	1.021070	1.018854	1.016501		
0.189475	1.030278	1.028910	1.027337	1.025586	1.023663	1.021581	1.019352	1.016986		
0.199415	1.030725	1.029345	1.027763	1.025998	1.024065	1.021770	1.019741	1.017367		

Table. 2 : Apparent molar volume ($\phi_v/\text{cm}^3\cdot\text{mol}^{-1}$) of SDS in aqueous and aqueous amino acid solutions as a function of molality and different temperature.

WATER+SDS m (mol.kg ⁻¹) (SDS), \sqrt{m}	ϕ_v (cm ³ .mol ⁻¹)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
0.000340	179.61	191.48	185.68	191.73	197.83	198.02	198.24	201.48		
0.000890	209.83	214.46	216.88	219.35	219.60	224.43	223.61	227.40		
0.001360	223.80	227.63	230.78	231.77	233.54	236.10	237.22	238.38		
0.014990	232.91	235.08	236.84	238.72	240.38	241.88	243.09	244.69		
0.029990	236.93	238.75	240.38	241.94	243.41	244.82	246.15	247.45		
0.043980	238.16	239.92	241.48	242.95	244.36	245.75	247.10	248.32		
0.058070	238.95	240.66	242.16	243.63	245.08	246.45	247.71	248.97		
0.071994	239.37	241.05	242.58	244.03	245.44	246.74	248.04	249.24		
0.085940	239.66	241.28	242.79	244.22	245.64	246.96	248.25	249.48		
0.100850	239.85	241.46	242.99	244.39	245.76	247.11	248.37	249.61		
0.114980	240.33	241.92	243.42	244.82	246.18	247.50	248.76	249.99		
0.128601	240.54	242.10	243.57	244.98	246.34	247.65	248.91	250.15		
0.142570	240.88	242.45	243.92	245.30	246.67	247.97	249.22	250.44		
0.157000	241.00	242.56	244.03	245.41	246.77	248.08	249.33	250.55		
0.171940	239.61	241.80	243.31	244.76	246.15	247.50	248.79	250.05		
0.186100	240.54	242.37	243.82	245.22	246.58	247.88	249.14	250.36		
0.198450	241.18	242.72	244.18	245.56	246.91	248.20	249.44	250.65		

WATER+GLYCINE(m=1.3)+SDS

m (mol.kg ⁻¹) (SDS) \sqrt{m}	ϕ_v (cm ³ .mol ⁻¹)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
0.001592	0.039898	337.04	338.81	338.30	340.20	341.58	340.06	340.95		
0.013725	0.117154	246.12	247.66	249.04	250.47	252.00	252.13	253.67		
0.027987	0.167293	245.37	246.86	248.16	249.57	250.78	251.79	252.73		
0.039750	0.199374	239.81	241.15	242.48	243.74	244.93	246.02	246.81		
0.053373	0.231027	242.40	243.73	244.99	246.28	247.45	248.60	249.72		
0.069147	0.262959	244.17	245.51	246.79	248.00	249.24	250.36	251.50		
0.085642	0.292646	242.91	244.25	245.54	246.78	247.99	249.12	250.29		
0.100063	0.316328	245.04	246.36	247.63	248.85	250.05	251.19	252.36		
0.114261	0.338025	245.13	246.43	247.70	248.96	250.14	251.29	252.46		
0.127985	0.357749	244.66	245.95	247.22	248.45	249.65	250.81	251.93		
0.140583	0.374945	244.95	246.23	247.49	248.74	249.92	251.07	252.22		
0.156884	0.396086	243.44	244.77	246.01	247.25	248.46	249.61	250.76		
0.171385	0.413986	245.11	246.40	247.67	248.89	250.09	251.25	252.40		
0.185656	0.430879	244.12	245.42	246.68	247.91	249.10	250.25	251.41		
0.195929	0.442639	245.29	246.59	247.83	249.07	250.27	251.41	252.57		

WATER+GLYCINE(m=0.883)+SDS

m (mol.kg ⁻¹) (SDS) \sqrt{m}	ϕ_v (cm ³ .mol ⁻¹)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
0.000680	810.22	810.43	805.31	797.58	792.85	808.12	813.74	802.47		
0.001544	499.40	501.49	500.04	503.04	501.83	503.23	504.74	504.47		
0.014193	263.35	264.72	265.87	267.42	268.34	269.65	271.48	272.75		
0.024522	247.03	248.54	249.74	251.30	252.48	253.81	255.47	255.94		
0.037557	247.44	248.85	250.12	251.57	252.79	254.07	255.32	256.35		
0.051069	246.74	248.17	249.45	250.84	252.06	253.29	254.54	255.60		
0.071936	243.75	245.20	246.60	247.87	249.16	250.41	251.55	252.62		
0.086225	243.95	245.33	246.65	247.99	249.22	250.46	251.69	252.79		
0.100026	244.30	245.69	247.00	248.32	249.57	250.80	252.03	253.15		
0.114767	243.51	244.87	246.21	247.53	248.77	250.01	251.21	252.35		
0.128215	243.00	244.38	245.72	247.05	248.28	249.52	250.73	251.87		
0.142658	243.07	244.45	245.77	247.09	248.33	249.56	250.76	251.91		
0.157116	243.48	244.87	246.18	247.49	248.73	249.95	251.16	252.30		
0.171290	243.49	244.88	246.16	247.46	248.71	250.52	251.73	252.87		
0.185205	243.54	244.88	246.22	247.61	248.84	250.08	251.62	252.77		
0.198478	243.68	245.06	246.36	247.67	248.90	250.13	251.34	252.48		

WATER+DL-ALANINE(1.216)+SDS

m (mol.kg ⁻¹) (SDS \sqrt{m})	ϕ_v (cm ³ .mol ⁻¹)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
0.000760	0.027568	712.65	709.65	710.14	710.83	710.36	730.08	732.69		
0.001556	0.039450	266.61	267.77	269.13	270.47	271.58	274.72	275.80		
0.014299	0.119578	241.37	244.17	245.60	246.93	248.16	250.29	251.56		
0.028638	0.169229	243.48	244.86	247.53	248.87	250.04	252.16	253.53		
0.042769	0.206806	243.15	244.58	247.23	248.56	249.75	251.47	252.62		
0.055985	0.236611	243.37	244.76	247.45	248.74	249.95	251.53	252.73		
0.070499	0.265516	242.61	244.02	245.37	246.67	249.17	250.72	251.89		
0.083252	0.288534	243.06	244.46	245.79	247.13	249.58	251.07	252.25		
0.099087	0.314781	245.74	247.12	248.43	249.73	252.21	253.66	254.72		
0.112982	0.336128	245.81	247.18	248.50	249.79	252.24	253.65	254.86		
0.127136	0.356562	243.74	245.12	246.44	247.73	250.22	251.63	252.81		
0.139751	0.373833	243.46	244.85	246.18	247.48	249.97	251.34	252.52		
0.154131	0.392596	244.28	245.64	246.94	248.23	250.66	252.03	253.21		
0.169950	0.412250	245.09	246.45	247.75	249.05	251.47	252.81	253.99		
0.189475	0.435287	246.14	247.47	248.77	250.05	252.46	253.80	254.99		
0.199415	0.446559	714.54	712.65	709.65	710.14	710.36	730.08	732.69		

WATER+DL-ALANINE(0.827)+SDS

m (mol.kg ⁻¹) (SDS) \sqrt{m}	ϕ_v (cm ³ .mol ⁻¹)									
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C		
0.001430	0.037810	186.43	187.82	189.23	189.39	189.57	194.89	195.13		
0.013983	0.118250	235.43	236.92	238.45	239.40	241.02	243.30	244.38		
0.026784	0.163659	244.91	246.47	247.87	249.26	250.28	252.43	253.46		
0.041951	0.204819	243.21	244.64	246.10	247.38	248.46	250.20	251.33		
0.055598	0.235791	244.71	246.10	247.47	248.75	249.93	251.48	252.58		
0.070241	0.265030	242.74	244.12	245.48	246.75	247.96	249.39	250.61		
0.082870	0.287871	242.44	243.77	245.17	246.44	247.61	249.05	250.24		
0.098097	0.313204	242.82	244.23	245.56	246.83	248.01	249.42	250.57		
0.108665	0.329644	243.47	244.87	246.20	247.48	248.69	250.03	251.19		
0.125733	0.354588	243.45	244.83	246.18	247.44	248.64	249.99	251.15		
0.140410	0.374714	245.70	247.06	248.38	249.64	250.85	252.17	253.33		
0.153293	0.391526	244.00	245.36	246.71	247.98	249.19	250.50	251.65		
0.168333	0.410284	246.31	247.65	249.07	250.43	251.42	252.72	253.95		
0.182099	0.426731	244.11	245.48	246.79	248.06	249.27	250.59	251.74		
0.198069	0.445049	244.81	246.18	247.48	248.75	249.96	251.26	252.42		

Table 3. Limiting apparent molar, ϕ_v° of SDS in aqueous and aqueous amino acid solutions as a function of molality and temperature. N= concentration number

System	N	m of SDS	$\phi_v^\circ / \text{cm}^3 \cdot \text{mol}^{-1}$							
			15°C	20°C	25°C	30°C	35°C	40°C	45°C	50°C
Water+SDS	17	0.00034- 0.19845	221.16	225.53	229.07	230.54	232.76	235.77	237.28	239.02
Water+SDS +										
Gly(0.883m)	13	0.02452- 0.19848	253.16	255.06	256.16	258.55	259.78	261.67	264.55	264.18
Water+SDS +										
Gly(1.3m)	14	0.01372- 0.19593	242.84	243.38	245.62	247.45	249.52	251.70	251.04	253.05
Water+SDS +										
DL-ala(0.827m)	15	0.00156- 0.19941	230.33	232.14	234.06	236.00	237.36	239.56	242.40	243.81
Water+SDS +										
DL-ala(1.216m)	13	0.02678- 0.19807	233.76	235.69	237.25	239.28	241.10	242.68	246.92	249.05

Table 4: Critical micelle concentration (c.m.c) of SDS in aqueous and aqueous amino acid solutions as a function of molality and different temperature.

Critical micelle concentration (c.m.c) of SDS								
	WATER+SDS							
	15°C	20°C	25°C	30°C	35°C	40°C	45°C	50°C
SDS	0.016607	0.016811	0.016400	0.016370	0.016410	0.016340	0.016330	0.016230
con.	WATER+GLYCINE(0.883m)+SDS							
	0.020989	0.020993	0.020990	0.021008	0.021003	0.020998	0.021023	0.021017
	WATER+GLYCINE(1.3m)+SDS							
(m/mol.kg ⁻¹)	0.024469	0.024452	0.024457	0.024524	0.024562	0.024600	0.024644	0.024734
(0.0-0.2m)	WATER+DL-ALANINE(0.827m)+SDS							
	0.023050	0.023000	0.022950	0.022887	0.022875	0.022827	0.022631	0.022691
	WATER+DL-ALANINE(1.216m)+SDS							
	0.032913	0.032919	0.032914	0.032919	0.032921	0.032919	0.032966	0.032977

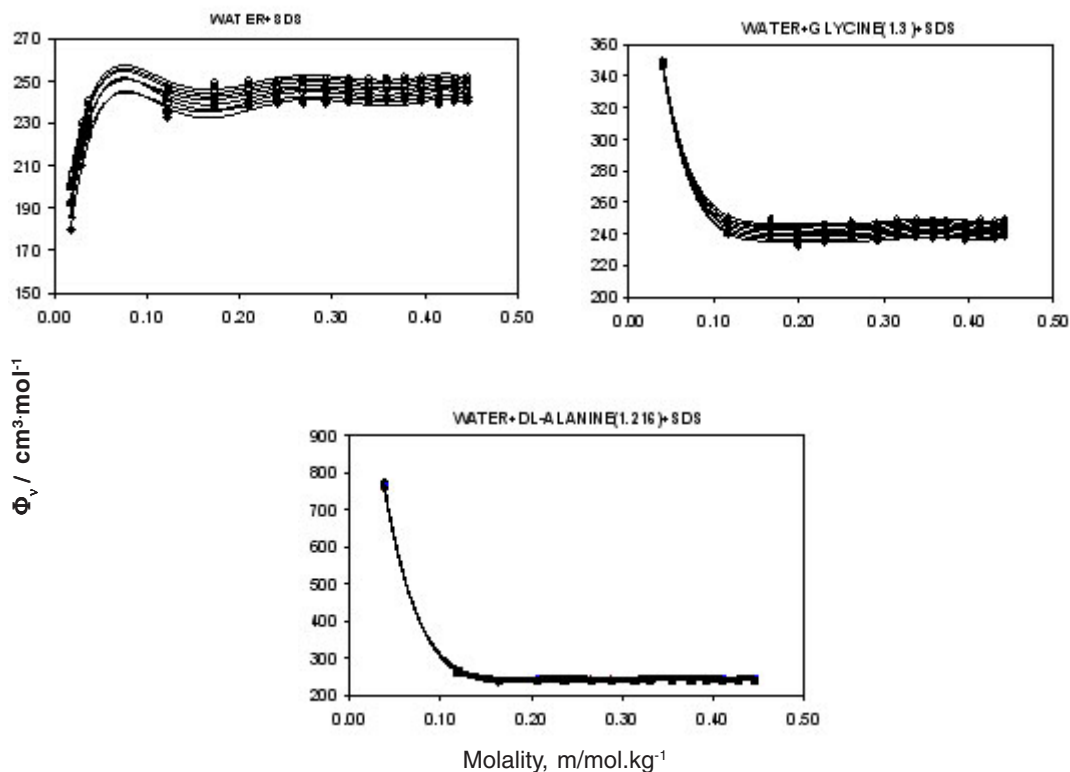


Fig 1: Limiting apparent molar volume of SDS in water , water+glycine(1.3), dl-alanine (1.216) as a function of molality of SDS at different temperature; ◆-15°C, ■-20 °C, ▲-25 °C, ◇-30 °C, □-35 °C, △-40 °C, ●-45 °C, ○-50 °C.

amino acid solutions are presented in Table 4.

The ϕ_v versus \sqrt{m} data were fitted to the equation

$$\phi_v / \text{cm}^3 \cdot \text{mol}^{-1} = \phi_v^0 + S_v m^{1/2} \quad \dots(2)$$

by least square method. The limiting apparent molar volumes, ϕ_v^0 , of SDS are presented in Table 3.

Table 4 reveals that our c.m.c values of SDS in water are somewhat larger than the literature values (≈ 0.009). An increase in c.m.c values in glycine and alanine solution was observed which demonstrates that glycine and alanine behave as water structure breakers. It has been reported that addition of structure breakers causes an increase of cmc¹.

The viscometric study of Devine and

Lowe² reports that glycine is water structure breaker whereas, alanine is water structure maker. The effect of concentration of glycine and alanine and that of temperature is not regular, the observed trend is that, c.m.c increases with the increase in amino acid concentration. c.m.c's of SDS would be expected to increase with increasing temperature¹. No such distinct behaviour is observed from the study. This may be due to the uncertainties associated with the measurement in very low concentration region. The uncertainty in measurement is greater than the expected variation of cmc values. Further attempts of cmc determination through ϕ_v measurement will establish this method.

It is evident from Figs. 1-3 that values of SDS increase with the increase in solute concentration. The concentration dependence of in aqueous and mixed aqueous solutions has been explained in terms of the solute-solute interaction. The usual interpretation is that the solute species

interact through the destructive overlap of their hydration spheres³. For apolar species, the positive volume contribution to ϕ_v , originated from the hydrophobic hydration starts decreasing as the solute concentration is increased. The overlap of two hydrophobic hydration cospheres relaxes some water molecules from the solvation sphere to the bulk giving rise to a negative change in volume. For hydrophobic ionic species the volume of water molecules is smaller in the solvation shell due to (a) the effect of electrostriction⁴ and (b) a decrease in the hydrogen-bonded network of water molecules in the solvation sphere than in the bulk (the so called structure-breaking effect). Further, if the ions are oppositely charged, this type of interaction also causes an attraction since the orientation of water molecules by the cations leads to favorable anion-water dipole interactions (and vice versa). In this way even the electrostricted water molecules may be shared, resulting in a positive volume change. The structure-breaking influence of ionic species on the hydrophobic hydration sphere of apolar groups gives a negative volume effect

Anionic surfactant sodium dodecyl sulphate has a long hydrophobic chain and charge centre. The concentration dependence ϕ_v of SDS can be interpreted in terms of the interactions involving the hydrophobic part as well as the charged centres. The ϕ_v values of SDS increase with the increase in molality i.e. S_v value positive. The positive sign of S_v indicates that interactions involving charged moieties dominate over the apolar group-apolar group and apolar group-charged centre interactions.

The ϕ_v^0 values, obtained by fitting $\phi_v - \sqrt{m}$ data by least squares method are tabulated in Table 3. The values of SDS in aqueous amino acid solutions are greater than these in water and increase with the concentration at all the studied temperatures. The variation of ϕ_v^0 with molality of amino acids can be rationalized in terms of the

cosphere overlap model⁵. According to this model, the overlap of cosphere of ion (or a polar group) with that of another ion (or a polar group) gives positive change in volume, while that of an ion (or a polar group) with that of a hydrophobic group and a hydrophilic group with another hydrophobic group results in a negative change in volume. As in the limit of infinite dilution the solute-solute interaction is minimal, the mutual overlap of the cospheres of zwitterionic amino acid with those of hydrophobic and hydrophilic group of SDS produces negative and positive changes in volume, respectively. The greater ϕ_v^0 values in amino acid solution demonstrate strong ion-ion interactions in the ternary system. The relaxation of the electrostricted water molecules, due to strong localized ion-zwitterion interactions, from the cospheres of both the amino acid and SDS to the bulk causes an increase in volume, this increase in volume outweighs the volume decrease due to the overlap of cospheres of hydrophobic parts of SDS and amino acids. Overlap of cosphere of amino acids gives positive change in volume.

As the concentration of amino acid increases the zwitterions-SDS anion and also the amino acid interactions increase giving rise to an increased ϕ_v^0 value. This in, and all probability account for the increase in ϕ_v^0 value with the increase in amino acid concentration.

The ϕ_v^0 of SDS in ternary solution can alternatively be thought of as arising from four constituents⁶, as

$$\phi_v^0 = V_{vw} + V_f + V_h + V_s \quad \dots(3)$$

where, V_{vw} and V_f are van der Waals volume⁷ and volume of empty spaces present therein⁸, respectively. The V_h and V_s represent the contributions due to the hydrophobic and hydrophilic hydration. The V_{vw} and V_f are assumed to be the same in aqueous amino acid as in water⁹.

Table 5: α_0 is the expansibility of water taken from G. S. Kell¹⁰

Temp.	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
α_0	1.51E-04	2.00E-04	2.57E-04	2.90E-04	3.46E-04	3.80E-04	4.22E-04	4.70E-04

Table 6: The parameters a_1 and a_2

w= Water;

sds= sodium dodecyl sulphate

w+sds		w+glycine(1.3)+sds+w+glycine(0.883)+sds						
m	a1	a2	m	a1	a2	m	a1	a2
0.000340	-7.81E-07	8.57E-09	0.001592	-7.44E-07	1.12E-08	0.00068	1.52E-06	-1.74E-08
0.000890	-6.85E-07	5.00E-09	0.014431	-2.44E-06	-9.52E-10	0.001544	3.98E-07	-2.38E-09
0.001360	-1.36E-06	1.33E-08	0.013725	-5.32E-06	4.26E-08	0.014193	-2.92E-06	4.29E-09
0.014990	-7.20E-06	4.93E-08	0.027987	-1.01E-05	7.55E-08	0.024522	-8.55E-06	5.79E-08
0.029990	-1.45E-05	1.18E-07	0.039750	-1.45E-05	1.11E-07	0.037557	-1.19E-05	7.55E-08
0.043980	-1.61E-05	9.88E-08	0.053373	-1.52E-05	8.64E-08	0.051069	-1.56E-05	9.64E-08
0.058070	-2.11E-05	1.31E-07	0.069147	-1.98E-05	1.14E-07	0.071936	-2.34E-05	1.56E-07
0.071994	-2.62E-05	1.69E-07	0.085642	-2.50E-05	1.46E-07	0.086225	-2.46E-05	1.37E-07
0.085940	-3.04E-05	1.90E-07	0.100063	-2.75E-05	1.53E-07	0.100026	-2.88E-05	1.63E-07
0.100850	-3.54E-05	2.22E-07	0.114261	-3.15E-05	1.76E-07	0.114767	-3.30E-05	1.88E-07
0.114980	-3.96E-05	2.48E-07	0.127985	-3.51E-05	1.97E-07	0.128215	-3.66E-05	2.05E-07
0.128601	-4.27E-05	2.57E-07	0.140583	-3.80E-05	2.10E-07	0.142658	-4.04E-05	2.26E-07
0.142570	-4.70E-05	2.87E-07	0.156884	-4.33E-05	2.42E-07	0.157116	-4.45E-05	2.52E-07
0.157000	-5.10E-05	3.06E-07	0.171385	-4.61E-05	2.55E-07	0.17129	-4.30E-05	1.40E-07
0.171940	-7.02E-05	5.09E-07	0.185656	-4.96E-05	2.71E-07	0.185205	-4.68E-05	1.80E-07
0.186100	-6.62E-05	4.40E-07	0.195929	-5.17E-05	2.81E-07	0.198478	-5.52E-05	3.11E-07
0.198450	-6.41E-05	3.92E-07						

w+dl-alanine(1.216)+sds			w+dl-alanine(.827)+sds		
m	a1	a2	m	a1	a2
0.001430	3.68E-06	-6.05E-08	0.000760	-3.85E-07	2.14E-09
0.013983	-2.02E-06	-1.86E-08	0.001556	-6.08E-07	3.10E-09
0.026784	-6.13E-06	-2.38E-10	0.014299	-3.66E-06	8.57E-09
0.041951	-7.61E-06	-2.41E-08	0.028638	-9.11E-06	4.74E-08
0.055598	-1.45E-05	4.74E-08	0.042769	-1.24E-05	5.95E-08
0.070241	-1.78E-05	5.95E-08	0.055985	-1.66E-05	8.33E-08
0.082870	-2.18E-05	8.36E-08	0.070499	-1.95E-05	8.86E-08
0.098097	-2.62E-05	1.11E-07	0.083252	-2.49E-05	1.34E-07
0.108665	-2.89E-05	1.33E-07	0.099087	-3.02E-05	1.71E-07
0.125733	-3.20E-05	1.33E-07	0.112982	-3.37E-05	1.86E-07
0.140410	-3.67E-05	1.59E-07	0.127136	-3.66E-05	1.99E-07
0.153293	-4.17E-05	1.99E-07	0.139751	-4.12E-05	2.29E-07
0.168333	-4.33E-05	1.95E-07	0.154131	-4.71E-05	2.86E-07
0.182099	-4.77E-05	2.28E-07	0.169950	-4.88E-05	2.64E-07
0.198069	-5.01E-05	2.31E-07	0.189475	-5.38E-05	2.92E-07
			0.199415	-6.84E-05	4.65E-07

Table 7: Expansibility (α) and apparent molar expansion (ϕ_E) of SDS in water as a function of molality and different temperature.

WATER+SDS								
m (mol.Kg⁻¹)	Expansibility ($\alpha \times 10^{-4}$)/C⁻¹							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.000340	1.5139	2.0043	2.5755	2.9026	3.4590	3.8008	4.2245	4.6991
0.000890	1.5139	2.0047	2.5763	2.9037	3.4604	3.8026	4.2266	4.7016
0.001360	1.5182	2.0081	2.5789	2.9054	3.4614	3.8027	4.2259	4.7000
0.014990	1.5646	2.0507	2.6175	2.9403	3.4924	3.8300	4.2493	4.7197
0.029990	1.6164	2.0954	2.6551	2.9710	3.5159	3.8465	4.2588	4.7220
0.043980	1.6365	2.1171	2.6784	2.9960	3.5426	3.8750	4.2890	4.7539
0.058070	1.6755	2.1527	2.7105	3.0248	3.5680	3.8971	4.3078	4.7693
0.071994	1.7140	2.1872	2.7410	3.0514	3.5906	3.9157	4.3225	4.7800
0.085940	1.7489	2.2197	2.7711	3.0793	3.6161	3.9391	4.3435	4.7988
0.100850	1.7875	2.2549	2.8028	3.1077	3.6410	3.9607	4.3617	4.8135
0.114980	1.8206	2.2852	2.8303	3.1326	3.6631	3.9801	4.3785	4.8276
0.128601	1.8474	2.3109	2.8548	3.1561	3.6855	4.0014	4.3987	4.8468
0.142570	1.8810	2.3413	2.8821	3.1802	3.7065	4.0194	4.4136	4.8584
0.157000	1.9134	2.3716	2.9102	3.2064	3.7304	4.0413	4.4334	4.8762
0.171940	2.0422	2.4802	2.9985	3.2744	3.7780	4.0686	4.4402	4.8624
0.186100	2.0219	2.4665	2.9913	3.2740	3.7843	4.0816	4.4601	4.8890
0.198450	2.0146	2.4636	2.9930	3.2803	3.7952	4.0972	4.4803	4.9140
	Apparent molar expansion (ϕ_E /cm³.mol⁻¹.C⁻¹)							
0.000340	0.02872	0.03965	0.04884	0.05641	0.06894	0.07552	0.08373	0.09441
0.000890	0.03236	0.04352	0.05635	0.06411	0.07635	0.08564	0.09475	0.10710
0.001360	0.03468	0.04631	0.06002	0.06774	0.08114	0.08999	0.10035	0.11204
0.014990	0.03681	0.04855	0.06230	0.07046	0.08419	0.09284	0.10347	0.11562
0.029990	0.03866	0.05035	0.06410	0.07212	0.08578	0.09433	0.10495	0.11692
0.043980	0.03927	0.05106	0.06492	0.07301	0.08676	0.09540	0.10613	0.11817
0.058070	0.04032	0.05207	0.06588	0.07391	0.08763	0.09621	0.10685	0.11886
0.071994	0.04131	0.05298	0.06673	0.07468	0.08831	0.09678	0.10735	0.11925
0.085940	0.04219	0.05381	0.06751	0.07541	0.08901	0.09744	0.10797	0.11983
0.100850	0.04315	0.05470	0.06833	0.07616	0.08967	0.09803	0.10847	0.12026
0.114980	0.04402	0.05553	0.06912	0.07690	0.09036	0.09866	0.10906	0.12080
0.128601	0.04470	0.05619	0.06976	0.07752	0.09097	0.09925	0.10963	0.12136
0.142570	0.04557	0.05701	0.07052	0.07821	0.09160	0.09982	0.11013	0.12179
0.157000	0.04637	0.05776	0.07123	0.07888	0.09223	0.10041	0.11067	0.12228
0.171940	0.04924	0.06025	0.07321	0.08036	0.09318	0.10085	0.11060	0.12168
0.186100	0.04891	0.06003	0.07316	0.08049	0.09349	0.10133	0.11124	0.12251
0.198450	0.04884	0.06003	0.07329	0.08074	0.09388	0.10184	0.11189	0.12328
WATER+GLYCINE(1.3)+SDS								
m (mol.Kg⁻¹)	Expansibility ($\alpha \times 10^{-4}$)/C⁻¹							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.001592	1.5129	2.0031	2.5653	2.9010	3.4573	3.7989	4.2223	4.6967

0.014431	1.5333	2.0237	2.5879	2.9249	3.4823	3.8251	4.2498	4.7255
0.013725	1.5484	2.0339	2.5940	2.9264	3.4793	3.8176	4.2377	4.7088
0.027987	1.5855	2.0661	2.6240	2.9531	3.5024	3.8374	4.2540	4.7217
0.039750	1.6177	2.0933	2.6484	2.9737	3.5192	3.8503	4.2632	4.7270
0.053373	1.6322	2.1094	2.6676	2.9953	3.5431	3.8766	4.2919	4.7581
0.069147	1.6684	2.1413	2.6979	3.0228	3.5676	3.8984	4.3108	4.7741
0.085642	1.7103	2.1779	2.7327	3.0541	3.5954	3.9228	4.3317	4.7916
0.100063	1.7327	2.1986	2.7536	3.0743	3.6149	3.9416	4.3498	4.8090
0.114261	1.7654	2.2274	2.7812	3.0994	3.6375	3.9617	4.3674	4.8241
0.127985	1.7941	2.2525	2.8052	3.1213	3.6570	3.9790	4.3825	4.8369
0.140583	1.8185	2.2745	2.8268	3.1415	3.6758	3.9965	4.3987	4.8517
0.156884	1.8602	2.3106	2.8612	3.1724	3.7031	4.0203	4.4189	4.8684
0.171385	1.8851	2.3331	2.8834	3.1934	3.7228	4.0387	4.4361	4.8842
0.185656	1.9138	2.3585	2.9083	3.2165	3.7439	4.0581	4.4536	4.8999
0.195929	1.9316	2.3746	2.9241	3.2313	3.7578	4.0710	4.4655	4.9109

Apparent molar expansion (Φ_E /cm³.mol⁻¹.C⁻¹)

0.001592	0.05108	0.06770	0.08650	0.09820	0.11761	0.12969	0.14345	0.15993
0.014431	0.04200	0.05568	0.07137	0.08109	0.09705	0.10715	0.11970	0.13394
0.013725	0.03819	0.05030	0.06440	0.07307	0.08730	0.09633	0.10694	0.11951
0.027987	0.03897	0.05093	0.06496	0.07347	0.08756	0.09636	0.10721	0.11941
0.039750	0.03882	0.05042	0.06405	0.07229	0.08593	0.09443	0.10498	0.11673
0.053373	0.03956	0.05133	0.06519	0.07355	0.08741	0.09607	0.10682	0.11892
0.069147	0.04073	0.05248	0.06641	0.07477	0.08863	0.09730	0.10805	0.12017
0.085642	0.04154	0.05310	0.06693	0.07517	0.08888	0.09742	0.10803	0.12003
0.100063	0.04245	0.05407	0.06801	0.07630	0.09011	0.09870	0.10938	0.12147
0.114261	0.04326	0.05479	0.06871	0.07694	0.09071	0.09924	0.10987	0.12189
0.127985	0.04387	0.05530	0.06917	0.07733	0.09101	0.09947	0.11004	0.12196
0.140583	0.04452	0.05590	0.06978	0.07792	0.09158	0.10002	0.11056	0.12247
0.156884	0.04525	0.05644	0.07021	0.07821	0.09171	0.10002	0.11042	0.12218
0.171385	0.04617	0.05737	0.07122	0.07926	0.09281	0.10114	0.11157	0.12338
0.185656	0.04667	0.05776	0.07155	0.07951	0.09296	0.10122	0.11157	0.12329
0.195929	0.04733	0.05843	0.07228	0.08024	0.09374	0.10202	0.11238	0.12414

WATER+GLYCINE(0.883)+SDS

m (mol.Kg ⁻¹)	Expansibility ($\alpha \times 10^{-4}$)/C ⁻¹							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.000680	1.4996	1.9927	2.5667	2.8964	3.4556	3.8001	4.2265	4.7039
0.001544	1.5060	1.9978	2.5703	2.8985	3.4562	3.7993	4.2242	4.7000
0.014193	1.5355	2.0263	2.5979	2.9254	3.4822	3.8245	4.2487	4.7237
0.024522	1.5737	2.0591	2.6252	2.9474	3.4987	3.8357	4.2544	4.7239
0.037557	1.6007	2.0843	2.6485	2.9689	3.5184	3.8536	4.2704	4.7382
0.051069	1.6298	2.1111	2.6732	2.9914	3.5387	3.8718	4.2865	4.7521
0.071936	1.6865	2.1617	2.7176	3.0299	3.5710	3.8981	4.3067	4.7662
0.086225	1.7024	2.1792	2.7367	3.0508	3.5935	3.9224	4.3327	4.7939
0.100026	1.7349	2.2091	2.7639	3.0753	3.6155	3.9417	4.3495	4.8080
0.114767	1.7677	2.2393	2.7914	3.1004	3.6379	3.9616	4.3668	4.8228
0.128215	1.7962	2.2659	2.8162	3.1234	3.6591	3.9811	4.3845	4.8387
0.142658	1.8259	2.2934	2.8414	3.1465	3.6799	3.9997	4.4010	4.8529
0.157116	1.8573	2.3222	2.8675	3.1700	3.7007	4.0180	4.4166	4.8659

0.171290	1.8746	2.3502	2.9062	3.2197	3.7613	4.0900	4.4997	4.9602
0.185205	1.8992	2.3707	2.9227	3.2322	3.7697	4.0940	4.4999	4.9563
0.198478	1.9412	2.4000	2.9391	3.2359	3.7605	4.0719	4.4647	4.9080

Apparent molar expansion (Φ_E /cm³.mol⁻¹.C⁻¹)

0.000680	0.12019	0.16045	0.20592	0.23049	0.27373	0.30711	0.34422	0.37803
0.001544	0.07504	0.10004	0.12841	0.14571	0.17338	0.19114	0.21319	0.23711
0.014193	0.04062	0.05382	0.06925	0.07841	0.09361	0.10330	0.11551	0.12901
0.024522	0.03913	0.05141	0.06577	0.07426	0.08850	0.09750	0.10881	0.12100
0.037557	0.03985	0.05209	0.06644	0.07487	0.08910	0.09805	0.10915	0.12156
0.051069	0.04045	0.05260	0.06688	0.07521	0.08935	0.09821	0.10923	0.12156
0.071936	0.04135	0.05323	0.06721	0.07528	0.08913	0.09775	0.10845	0.12049
0.086225	0.04175	0.05367	0.06769	0.07583	0.08971	0.09838	0.10917	0.12129
0.100026	0.04260	0.05448	0.06846	0.07654	0.09039	0.09900	0.10974	0.12182
0.114767	0.04327	0.05504	0.06891	0.07691	0.09065	0.09918	0.10982	0.12181
0.128215	0.04386	0.05558	0.06938	0.07733	0.09100	0.09948	0.11006	0.12198
0.142658	0.04460	0.05626	0.07002	0.07792	0.09153	0.09996	0.11048	0.12235
0.157116	0.04544	0.05706	0.07077	0.07862	0.09220	0.10056	0.11105	0.12287
0.171290	0.04585	0.05775	0.07173	0.07986	0.09372	0.10263	0.11343	0.12558
0.185205	0.04646	0.05825	0.07215	0.08021	0.09397	0.10254	0.11337	0.12541
0.198478	0.04752	0.05901	0.07259	0.08031	0.09375	0.10199	0.11233	0.12402

WATER+DL-ALANINE(1.216)+SDS

m (mol.kg ⁻¹)	Expansibility ($\alpha \times 10^{-4}$)C ⁻¹							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.001430	1.4916	1.9891	2.5674	2.9013	3.4649	3.8137	4.2445	4.7263
0.013983	1.5333	2.0264	2.6003	2.9300	3.4891	3.8337	4.2601	4.7375
0.026784	1.5664	2.0574	2.6291	2.9569	3.5138	3.8564	4.2809	4.7562
0.041951	1.5869	2.0800	2.6538	2.9838	3.5429	3.8878	4.3146	4.7922
0.055598	1.6318	2.1178	2.6845	3.0075	3.5596	3.8974	4.3171	4.7875
0.070241	1.6598	2.1444	2.7098	3.0316	3.5823	3.9188	4.3372	4.8064
0.082870	1.6899	2.1720	2.7348	3.0542	3.6024	3.9366	4.3525	4.8192
0.098097	1.7236	2.2030	2.7631	3.0798	3.6252	3.9567	4.3699	4.8338
0.108665	1.7434	2.2206	2.7785	3.0930	3.6362	3.9655	4.3765	4.8382
0.125733	1.7719	2.2490	2.8066	3.1211	3.6642	3.9935	4.4044	4.8661
0.140410	1.8089	2.2832	2.8380	3.1499	3.6902	4.0168	4.4251	4.8840
0.153293	1.8443	2.3147	2.8655	3.1735	3.7098	4.0325	4.4368	4.8918
0.168333	1.8604	2.3309	2.8820	3.1902	3.7267	4.0498	4.4544	4.9096
0.182099	1.8922	2.3595	2.9073	3.2124	3.7457	4.0656	4.4669	4.9189
0.198069	1.9144	2.3813	2.9288	3.2336	3.7666	4.0861	4.4872	4.9389

Apparent molar expansion (Φ_E /cm³.mol⁻¹.C⁻¹)

0.001430	0.10543	0.14101	0.18187	0.20613	0.24681	0.27184	0.31125	0.34809
0.013983	0.04081	0.05421	0.06982	0.07906	0.09459	0.10435	0.11728	0.13092
0.026784	0.03802	0.05017	0.06440	0.07283	0.08697	0.09591	0.10735	0.11985
0.041951	0.03882	0.05111	0.06552	0.07405	0.08837	0.09741	0.10901	0.12171
0.055598	0.03989	0.05200	0.06621	0.07454	0.08866	0.09751	0.10872	0.12110
0.070241	0.04060	0.05269	0.06688	0.07520	0.08928	0.09812	0.10925	0.12162
0.082870	0.04121	0.05320	0.06730	0.07552	0.08950	0.09825	0.10928	0.12153

0.098097	0.04211	0.05406	0.06810	0.07629	0.09021	0.09891	0.10986	0.12207
0.108665	0.04305	0.05507	0.06921	0.07741	0.09142	0.10016	0.11115	0.12337
0.125733	0.04376	0.05578	0.06992	0.07813	0.09214	0.10088	0.11186	0.12414
0.140410	0.04430	0.05616	0.07012	0.07821	0.09205	0.10066	0.11149	0.12360
0.153293	0.04511	0.05687	0.07073	0.07871	0.09244	0.10095	0.11165	0.12365
0.168333	0.04565	0.05745	0.07135	0.07936	0.09313	0.10166	0.11240	0.12444
0.182099	0.04658	0.05834	0.07221	0.08017	0.09390	0.10238	0.11306	0.12506
0.198069	0.04732	0.05912	0.07303	0.08102	0.09480	0.10330	0.11402	0.12605

WATER+DL-ALANINE(0.827)+SDS

m (mol.kg ⁻¹)	Expansibility ($\alpha \times 10^{-4}$)C ⁻¹							
	15 °C	20 °C	25 °C	30 °C	35 °C	40 °C	45 °C	50 °C
0.000760	1.5117	2.0028	2.5746	2.9023	3.4594	3.8018	4.2261	4.7014
0.001556	1.5136	2.0046	2.5764	2.9039	3.4609	3.8033	4.2275	4.7027
0.014299	1.5411	2.0314	2.6024	2.9294	3.4857	3.8274	4.2511	4.7256
0.028638	1.5820	2.0683	2.6353	2.9584	3.5106	3.8485	4.2681	4.7386
0.042769	1.6093	2.0942	2.6598	2.9816	3.5325	3.8691	4.2874	4.7566
0.055985	1.6425	2.1248	2.6878	3.0072	3.5555	3.8897	4.3055	4.7723
0.070499	1.6685	2.1501	2.7123	3.0311	3.5787	3.9123	4.3275	4.7935
0.083252	1.7067	2.1838	2.7414	3.0557	3.5986	3.9277	4.3383	4.7997
0.099087	1.7470	2.2203	2.7742	3.0848	3.6240	3.9494	4.3564	4.8141
0.112982	1.7754	2.2470	2.7992	3.1083	3.6459	3.9697	4.3751	4.8312
0.127136	1.7994	2.2697	2.8206	3.1283	3.6646	3.9871	4.3912	4.8459
0.139751	1.8342	2.3013	2.8489	3.1536	3.6867	4.0061	4.4071	4.8586
0.154131	1.8741	2.3356	2.8776	3.1768	3.7043	4.0180	4.4133	4.8593
0.169950	1.8953	2.3586	2.9024	3.2035	3.7327	4.0486	4.4459	4.8938
0.189475	1.9351	2.3956	2.9364	3.2348	3.7612	4.0743	4.4688	4.9139
0.199415	2.0261	2.4697	2.9937	3.2752	3.7846	4.0816	4.4582	4.8861

Apparent molar expansion (Φ_E /cm ³ .mol ⁻¹ .C ⁻¹)								
0.000760	0.02837	0.03770	0.04868	0.05522	0.06578	0.07231	0.08257	0.09191
0.001556	0.03573	0.04748	0.06131	0.06949	0.08308	0.09188	0.10305	0.11509
0.014299	0.03773	0.04997	0.06435	0.07281	0.08708	0.09598	0.10749	0.11995
0.028638	0.03848	0.05054	0.06468	0.07301	0.08703	0.09579	0.10694	0.11923
0.042769	0.03937	0.05146	0.06566	0.07397	0.08804	0.09686	0.10796	0.12028
0.055985	0.03986	0.05180	0.06582	0.07401	0.08790	0.09660	0.10752	0.11973
0.070499	0.04044	0.05233	0.06631	0.07450	0.08836	0.09703	0.10792	0.12009
0.083252	0.04143	0.05324	0.06715	0.07522	0.08899	0.09756	0.10834	0.12038
0.099087	0.04252	0.05428	0.06813	0.07613	0.08985	0.09836	0.10906	0.12104
0.112982	0.04320	0.05492	0.06873	0.07670	0.09038	0.09885	0.10950	0.12145
0.127136	0.04418	0.05597	0.06988	0.07788	0.09164	0.10016	0.11086	0.12287
0.139751	0.04472	0.05636	0.07010	0.07798	0.09158	0.09997	0.11053	0.12238
0.154131	0.04613	0.05774	0.07146	0.07930	0.09293	0.10116	0.11165	0.12350
0.169950	0.04622	0.05778	0.07144	0.07924	0.09275	0.10107	0.11154	0.12331
0.189475	0.04732	0.05885	0.07248	0.08023	0.09372	0.10199	0.11241	0.12415
0.199415	0.04949	0.06059	0.07378	0.08111	0.09415	0.10243	0.11195	0.12323

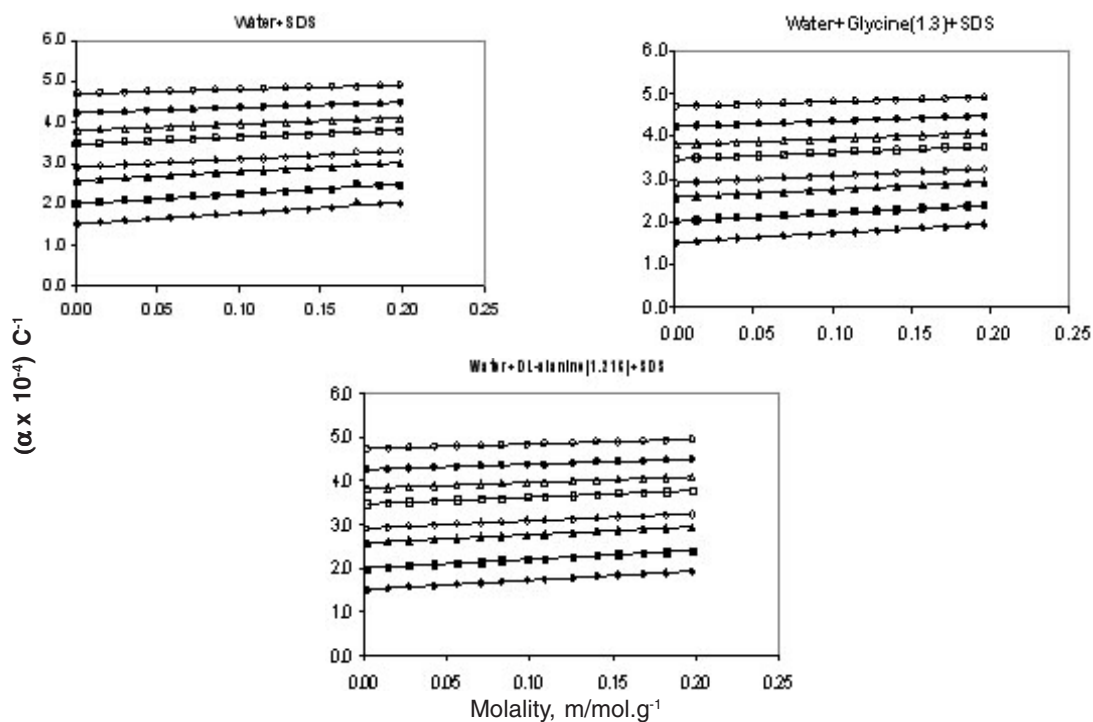


Fig 2:Expansibility (α) of SDS in water, water+glycine(1.3), water+dl-alanine (1.216) as a function of molality of SDS at different temperature; \blacklozenge -15°C, \blacksquare -20°C, \blacktriangle -25°C, \blacklozenge -30°C, \square -35°C, \triangle -40°C, \bullet -45°C, \circ -50°C.

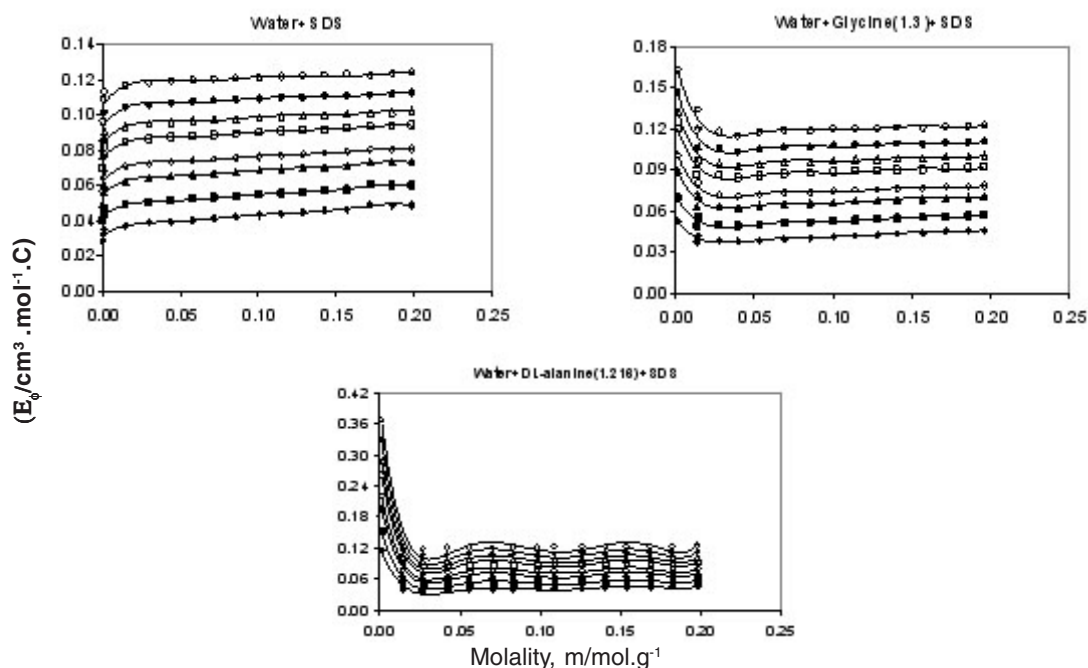


Fig. 3 Apparent molar expansion (ϕ_E) of SDS in water, water+glycine(1.3), water+dl-alanine(1.216) as a function of molality of SDS at different temperature; \blacklozenge -15°C, \blacksquare -20°C, \blacktriangle -25°C, \blacklozenge -30°C, \square -35°C, \triangle -40°C, \bullet -45°C, \circ -50°C.

The variation of ϕ_v^0 is, Therefore, due to the changes in $(V_n + V_s)$ resulting from SDS-amino acid, amino acid-amino acids and amino acid water interactions; the contribution from water-water interactions is assumed to be negligible.

At the same amino acid concentration, the ϕ_v^0 increases with the increase in temperature. This may be due to the resultant of the following effects;

- (i) Due to the increased thermal energy at higher temperature, the relaxation to the bulk of the electrostricted water molecules from the ionic interaction regions results in a positive volume change.
- (ii) An increase in temperature renders the amino acid- amino acid interaction relatively weaker giving rise to a small negative volume change.
- (iii) A decrease in amino acid-water interactions causes an increased volume change.

Apparent molar expansion and expansibility

Apparent molar expansion, ϕ_E was computed from measured density data using the relation

$$\phi_E = \frac{(\alpha - \alpha_0)}{m \rho_0} + \alpha \phi_v \quad (4)$$

where, α_0 is the expansibility of water taken from G. S. Kell¹⁰, α is the expansibility of solution calculated using the equation

$$\alpha = -\frac{a_1 + 2a_2T}{\rho} + \frac{\alpha_0 \rho_0}{\rho} \quad (5)$$

The parameters a_1 and a_2 were obtained by the regression of $(\rho - \rho_0) - T$ data according to equation

$$(\rho - \rho_0) / g.cm^{-3} = \sum_{i=0}^2 a_i T^i \quad (6)$$

using the a_1 parameter the expansibility of

solution was calculated from equation (5) and hence apparent molar expansion was calculated using equation (4). The calculated α and ϕ_E as a function SDS molality and temperature are presented in Table 7. Graphical representation of α and ϕ_E against molality is given in Fig 2-3. Apparent molar expansions in aqueous solution are very sensitive to hydrophobic and hydrophilic solute-solute interactions^{11,12}. Expansibility, and apparent molar expansion, ϕ_E reflect from the same structural alteration in the solutions. One might, therefore, expect the same response due to change of other parameters like concentration, temperature, pressure, etc. Hence, we concern with apparent molar expansion only. It is observed from isotherms that $\partial \phi_E / \partial m = a_1$ is positive for all the studied aqueous SDS and increases with an increase in the size of hydrophobic side chain of SDS and temperature. Since the aggregates due to hydrophobic interactions in aqueous environment are more sensitive to temperature and more unstable than the aggregates of water-water around the hydrophobic moieties, in turns, than that of around hydrophilic centers and in normal water¹⁶, the greater expansibility of the former would be expected.

The positive values of $\partial \phi_E / \partial m$ reveal that as an increase of molal concentration of SDS, the aggregation through hydrophobic moieties due to hydrophobic interactions are facile. The hydrophobic interactions are more dominant as the size of hydrophobic chain becomes larger¹⁶⁻¹⁸. The increase of $\partial \phi_E / \partial m$ values with the size of hydrophobic chain accounts the predominance of hydrophobic interactions in SDS with larger hydrocarbon backbone.

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