



RESEARCH ARTICLE

Antifungal Activities and Characterization of Some New Environmentally Safe Cu (II) Surfactants Substituted 2-Amino-6-Methyl Benzothiazole

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Abstract:**Background:**

Biologically potent compounds are one of the most important classes of materials for the upcoming generations.

Objectives:

Increasing number of microbial infectious diseases and resistant pathogens create a demand and urgency to develop novel, potent, safe and improved variety of antimicrobial agents.

Methods:

The copper surfactants substituted 2-amino-6-methyl benzothiazole were synthesized. The synthesized complexes have been characterized by IR, NMR, ESR spectroscopic methods. The antifungal activities have been evaluated by testing against *Alternaria alternate* fungi. All complexes showed good antifungal activity because chelation increases the anti-microbial potency.

Result:

The studies suggest that the copper (II) ions in soaps may be responsible for the enhancement of the activity against fungi. The evaluation of anti-fungal studies further revealed that fungitoxicity of the complexes also depends on the nature of metal ions. The chelation reduces the polarity of central metal ion mainly because of partial attaining of its positive charge with the donor groups and possible π -electron delocalization over the whole chelate ring. Such chelation increases the lipophilic character of the central atom, which subsequently favors its permeation through the lipid layer of the cell membrane. Their efficiency increases with their concentration.

Keywords: IR, Copper surfactants, Antifungal, *Alternaria alternate*, 2-amino-6-methyl benzothiazole, ESR spectroscopic method.

1. INTRODUCTION

Colloidal surfactants have become one of the most important requirements in day to day life, and have great importance in modern engineering and pharmacological fields. Many copper thiosemicarbazone complexes have been found to have significant anti-tubercular, fungicidal and antitumor activities [1, 2]. The use of copper linoleate as heavy duty wood preservative and many other biological activities of copper metal-containing surfactants have also been studied [3, 4]. Recently, the deeper understanding of the role of metal ion in bio-system has led to the awareness that metal complexing is useful in the treatment of bacterial, fungicidal and viral diseases [5, 6]. Many copper compounds are employed as pesticides either alone (as Cu (I) oxide and Cu (II) sulfate) or in mixtures [7]. Formulated proprietary

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brands of cuprous oxide are extensively employed as fungicides and seed dressings. Copper oxychloride has a number of applications, by far the most important being as an agricultural fungicide for which purpose it is extensively employed in formulated form as dust, wettable powders and pastes [8 - 10]. The effect of silver nanoparticles on various fungi has been reported recently [11, 12]. Therefore, a substantial research is needed for their discovery and improvement. Chemistry of present era aims to build a pollution free environment. For the same, it targets to create some alternatives which are eco-friendly and nature loving. The present research work is a step towards achieving such alternatives. For this, the copper surfactants substituted 2-amino-6-chloro benzothiazole were synthesized. The antifungal activities of these have been evaluated by testing against *Alternaria alternata*. The fungitoxicity results indicate that the strain of fungal species are susceptible towards these and suggests that with the increase in the concentration of copper surfactant and its complexes, it may increase further.

2. EXPERIMENTAL

The acids (n-caprylic acid, n-capric acid and n-lauric acid) have been purified by keeping over anhydrous sodium sulphate for a week and then distilled under reduced pressure. The purity of acids was checked by the determination of their melting points. n - Caprylic acid (16.3 °C) n - Capric acid (31.2 °C), n - Lauric acid (44.0 °C). Copper soaps were prepared by direct metathesis of the corresponding sodium soap with a slight excess of the required amount of copper sulphate solution at 50-55°C. After washing with hot water and alcohol, the samples were dried at 100-105°C. Finally, under reduced pressure, they were recrystallized twice from hot benzene. Ligands were synthesized using thiocyanation method [13]. The purity of the benzothiazoles has been checked by thin layer chromatography in various non-aqueous solvent systems. The metal was analysed by standard procedure. C, H, N, Cl and S analyses were performed at RSIC, CDRI Lucknow. In general, all the solid complexes (about 90% yield) with bluish green periphery were obtained (Table 1). These are soluble in ethanol, methanol, benzene and other organic solvents and insoluble in water. All the complexes are quite stable at room temperature up to 170°C. On the basis of elemental analysis, the complexes have been assigned the composition $Cu_2(C_nH_{2n+1}COO)_4L_2$, n=7,9,11 and suggested 1:1 type stoichiometry (Figs. (1-3) [14].

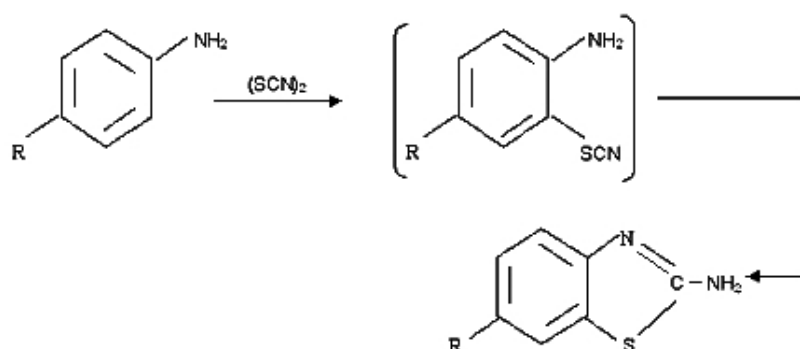


Fig. (1). Synthesis of ligand by thiocyanation method.

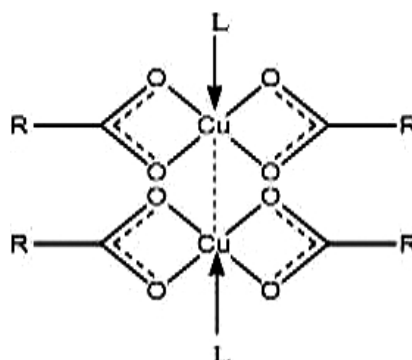


Fig. (2). Proposed structure of copper soap.

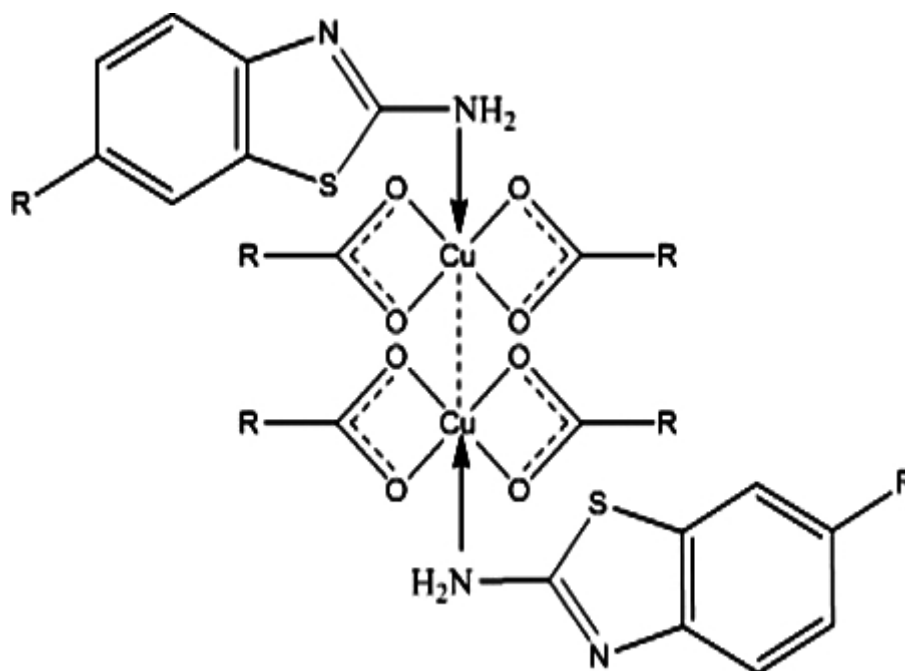


Fig. (3). Proposed structure of copper soap complex.

2.1. Fungicidal Activities:

1 ml of the soap solution was aseptically transferred into sterile petri plates. Into these plates, 20 ml of P.D.A. was poured and was mixed with soap solution by rotating the petri plates in clockwise and anticlockwise direction 3-4 times and was allowed to solidify. After the solidification of the above medium, single hypha / spore of test organism was aseptically transferred in the centre of the petri plates. The plates were incubated at 30 °C for 3 days. After the period of incubation, the plates were observed for the growth of fungus in different concentration of the soap solution used in the present study. The data were statistically analyzed according to the following formula [15].

$$\% \text{ Inhibition} = (C-T) \times 100 / C$$

Where C = diameter of fungal colony in control plates after 3 days

T = diameter of fungal colony in test plate after 3 days

3. CHARACTERIZATION

In order to study the structure of complexes, the infrared absorption spectra of compounds were obtained on a Perkin-Elmer 5100.4367 Spectrophotometer [4000-200 cm^{-1}] and ESR analysis was made on Marker-TCNE analyser at RSIC, IIT Powai, Mumbai. ^1H NMR spectra (C_6D_6) was recorded on an NMR spectrometer at RSIC, IIT Powai, Mumbai. The ESR spectra of the complex recorded on X-band at frequency 9.07 GHz under the magnetic field strength of 2000 ± 4000 and 3000 ± 2000 gauss at room temperature. ESR Spectrum is characterized by the position intensity and shape as to component lines. The position of ESR is referred as g values and is directly determined by the energy levels. The variation in g value is interpreted in terms of the first and second order spin-orbit interaction. The g values for ESR signal was calculated by the formulae [16].

$$g = hv / \beta H$$

Where,

ν = Freq. of band in kHz , β = Bohr magneton

H = magnetic field

h = Planck's constant

4. RESULTS AND DISCUSSION

4.1. IR Studies

The detailed infrared spectral investigation reveals that there are marked differences between the spectra of pure copper soaps and pure ligand than those of the corresponding complexes (Table 2) The absorption bands observed in the region $2960-2950\text{ cm}^{-1}$ and $2900-2950\text{ cm}^{-1}$ respectively were assigned to the antisymmetric stretching for methyl group of the soap segment present in the complex. Symmetric stretching bands for methyl group has been assigned in the region $2820-2850\text{ cm}^{-1}$. It has been observed that the absorption peaks, which are characteristics of the aliphatic portion of the soap molecules, remain unchanged on going while being transferred from soaps to the complexes. The absorption spectra in the region $710-725\text{ cm}^{-1}$ may probably be due to the rocking vibration of a chain of methylene group $-(\text{CH}_2)_n$. The peaks in the region $1575-1590\text{ cm}^{-1}$ and $1400-1410\text{ cm}^{-1}$ are due to COO^- , C-O anti-symmetric and symmetric stretching, respectively [17].

The presence of bands in the region $810\pm 20\text{ cm}^{-1}$ was due to C-H deformation (in plane) present in the tri-substituted benzene ring of benzothiazole segment, where two hydrogens are adjacent [1,2, and 4 position]. In 2-amino-6-methyl benzothiazole, the C-H deformation vibration of $-\text{CH}_3$ group has been observed at 1477 cm^{-1} . C-N stretching (aryl tertiary) peaks was observed in the region $1360-1310\text{ cm}^{-1}$. C=N stretching peaks in the region $1460-1430\text{ cm}^{-1}$ was also observed. The C-S stretching bands present in the region of $700-600\text{ cm}^{-1}$ also confirm the structure of the molecule. IR spectrum of the pure ligand 2-amino-6-R benzothiazole [R= $-\text{CH}_3$] contained peaks in the region $3450\pm 15\text{ cm}^{-1}$ due to symmetric N-H stretching, which shifted to $3400\pm 20\text{ cm}^{-1}$ on complexation, indicating the participation of the primary amino group in the chelation (Table 3) [18].

4.2. NMR Studies

The ^1H NMR spectra of the complexes were collected in C_6D_6 and examined to determine bonding and tautomerism present in the structure. A perusal of the spectra of the complexes shows aliphatic $-\text{CH}_3$ proton signal at $\delta - 0.9$ and $-\text{CH}_2$ proton signal at $\delta - 1.2$ of soap segment, indicating the presence of soap segment in the complex molecule.

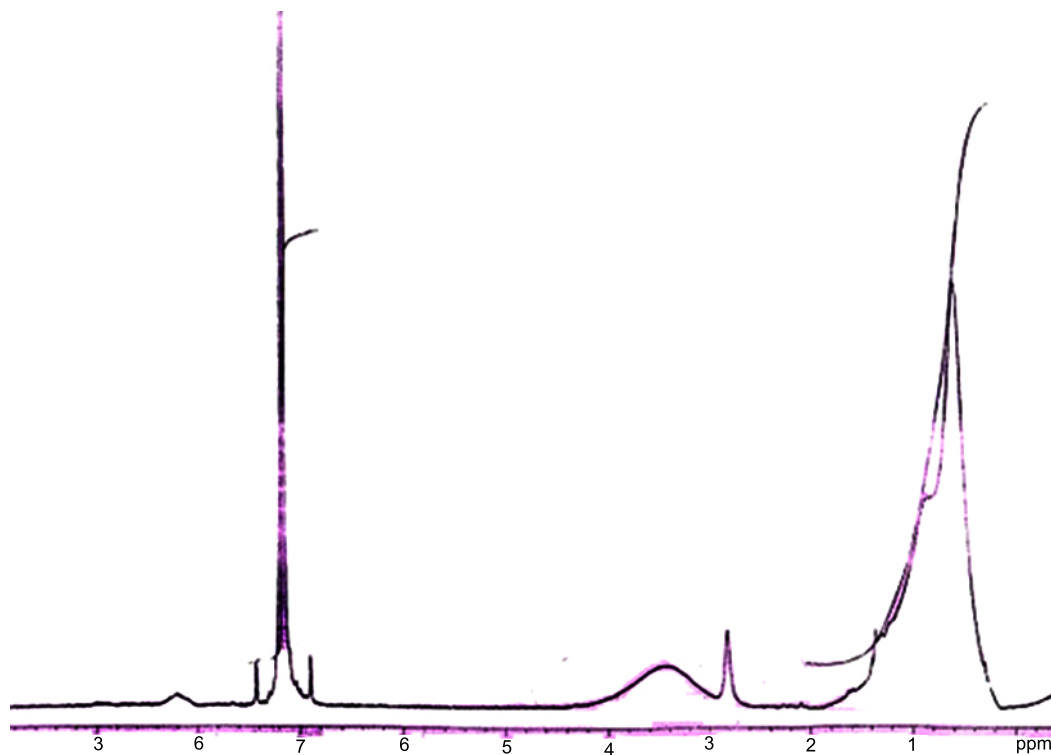


Fig. (4). NMR spectra of ligand (2-amino-6-methyl benzothiazole).

A broadened peak has been observed in all the spectra of the complexes at δ 3.2 – 4.2 due to the presence of aromatic $-\text{NH}_2$ protons. In the given spectra of $\text{CC}_{\text{pr}}\text{AMB}$ (Fig. 4), the above peak was clear which indicates the coordination through the $-\text{NH}_2$ group of the benzothiazole segment to the metal atom of soap segment. The proton present in the nitrogen atom may undergo rapid, intermediate or slow exchange. The broadening of the observed peak is suggestive of a slow exchange because the electrical quadrupole moment of the nitrogen nucleus induces a moderately efficient spin relaxation. A very weak signal was observed at δ - 8.5 in the spectra, which may be due to tautomerism present in the molecule of the complex similar to the pyrrole and indole. The triplets at δ 6.9 – 7.2 are may be due to tri-substituted aromatic protons of benzothiazole segment indicating that the environment of all the three protons is different. Another peak present at δ - 2.8 in the methyl substituted complex confirmed the presence of aromatic methyl group in the benzothiazole segment of the molecule suggesting a downfield shift [19].

4.3. ESR Studies

Literature reveals that magnetic conductivity and spectral studies of complexes of some first-row transition elements with pyrazine carboxylic acid have been conducted, where it acts as a bidentate ligand through O and N donor atoms [20].

Table 1. Analytical and physical data of the complexes.

Molecular Formula of Complex	Colour	Yield	M.P.	Cu	C	H	O	N	S
$\text{Cu}_2(\text{C}_7\text{H}_{13}\text{COO})_4 \cdot (\text{C}_8\text{H}_8\text{N}_2\text{S})_2$ $\text{CC}_{\text{pr}}\text{AMB}$	Dark Greenish Blue	92%	210	12.28 [12.36]	55.92 [56.09]	7.26 [7.40]	12.39 [12.45]	5.37 [5.45]	6.19 [6.23]
$\text{Cu}_2(\text{C}_9\text{H}_{19}\text{COO})_4 \cdot (\text{C}_8\text{H}_8\text{N}_2\text{S})_2$ $\text{CC}_{\text{pr}}\text{AMB}$	Greenish Blue	89%	220	10.80 [11.15]	58.87 [58.99]	8.01 [8.08]	11.15 [11.24]	4.80 [4.92]	5.52 [5.62]
$\text{Cu}_2(\text{C}_{11}\text{H}_{23}\text{COO})_4 \cdot (\text{C}_8\text{H}_8\text{N}_2\text{S})_2$ $\text{CL}_{\text{rt}}\text{AMB}$	Blue	91%	225	9.95 [10.15]	61.18 [61.39]	8.59 [8.63]	10.15 [10.23]	4.34 [4.47]	5.03 [5.11]

Table 2. Infrared absorption spectral frequencies (cm^{-1}) of soap and ligands and their assignments.

Assignment	Soap			Ligand
	Cu-Caprylate	Cu-Caprinate	Cu-Laurate	2-amino-6-methyl benzothiazole
N-H stretching in amino group	---	---	---	3434.9
CH_3 , C-H antisym. Stretching	2900.0	2950.0	2965.0	---
CH_2 , C-H symmetric stretching	2820.0	2840.0	2855.0	---
COO^- , C-O antisym. Stretching	1575.0	1590.0	1605.0	---
C-H deformation vibration of $-\text{CH}_3$ group	---	---	---	1477.9
C=N stretching	---	---	---	1460.0
CH_2 , deformation	1430.0	1430.0	1430.0	---
COO^- , C-O symmetric	1410.0	1410.0	1410.0	---
C-N stretching (aryl tertiary)	---	---	---	1350.0
CH_2 twisting and wagging	1295.0	1275.0 1230.0	1260.0	---
CH_3 rocking	1105.0	1110.0	1115.0	---
C-H deformation (in plane) due to benzene ring	---	---	---	808.1
CH_2 rocking	710.0	718.0	726.0	---
C-S stretching	---	---	---	680.0

Table 3. Infrared absorption spectral frequencies (cm^{-1}) of complexes and their assignments.

Assignment	$\text{CC}_{\text{pr}}\text{AMB}$	$\text{CC}_{\text{pr}}\text{AMB}$	$\text{CL}_{\text{rt}}\text{AMB}$
N-H stretching in amino group	3440.1	3419.2	3419.2
CH_3 , C-H antisym. Stretching	2932.6	2958.7	2922.1
CH_2 , C-H symmetric stretching	2864.5	2864.5	2848.8
COO^- , C-O antisym. Stretching	1590.0	1560.0	1575.0
C-H deformation vibration of $-\text{CH}_3$ group	1515.0	1520.0	1470.0

(Table 3) contd....

Assignment	CC _{pl} AMB	CC _{pr} AMB	CL _{rt} AMB
C=N stretching	1425.6	1420.3	1415.1
C-N stretching (aryl tertiary)	1341.9	1336.6	1350.0
CH ₃ rocking	1116.9	1111.6	1100.0
C-H deformation (in plane) due to benzene ring	829.1	815.0	805.0
CH ₂ rocking	705.0	740.0	724.4
C-S stretching	682.6	672.1	635.5

The ESR spectra of the complexes synthesized by us were recorded on X-band at frequency 9.07 GHz under the magnetic field strength of 2000±4000 and 3000±2000 gauss at room temperature. The value of ESR parameters for all the complexes is given in Table 4. A perusal of Table 4 shows that the values of g , $g_{||}$, g_{\perp} are greater than the value of g i.e. 2.0027. This indicates that the distortion from the regular octahedron has taken place in the shape of the complex. Also the trend $g_{||} > g_{\perp}$ for all the complexes indicates that the unpaired electron is most likely in $d_{x^2-y^2}$ orbital of Cu (II) giving $^2B_{1g}$ as the ground state. This fact supports that complexes possess elongated octahedral geometry. It is well known that $g_{||}$ is a moderately sensitive function for indicating covalency. Thus, the values of $g_{||}$ are suggestive of the covalent character of metal-ligand bond [21, 22]. Earlier magnetic studies, about copper (II) soaps confirm the binuclear configuration in the solid state. However, the low values of the magnetic moments for copper ion present in copper (II) soaps in non-aqueous solvent like alcohol and hydrocarbon confirm that the binuclear configuration exists in the soap solutions and the dispersion unit is a binuclear molecule having the structure proposed in Figs. (5-7) [23, 24].

Table 4. ESR spectral values of complexes.

Name of Complex and Molecular Formula	g	$g_{ }$	g_{\perp}
CC _{pl} AMB $Cu_2(C_7H_{15}COO)_4(C_8H_8N_2S)_2$	2.0184	2.2282	2.1282
CC _{pr} AMB $Cu_2(C_9H_{19}COO)_4(C_8H_8N_2S)_2$	2.0195	2.2349	2.1090
CL _{rt} AMB $Cu_2(C_{11}H_{23}COO)_4(C_8H_8N_2S)_2$	2.0270	2.2202	2.1092

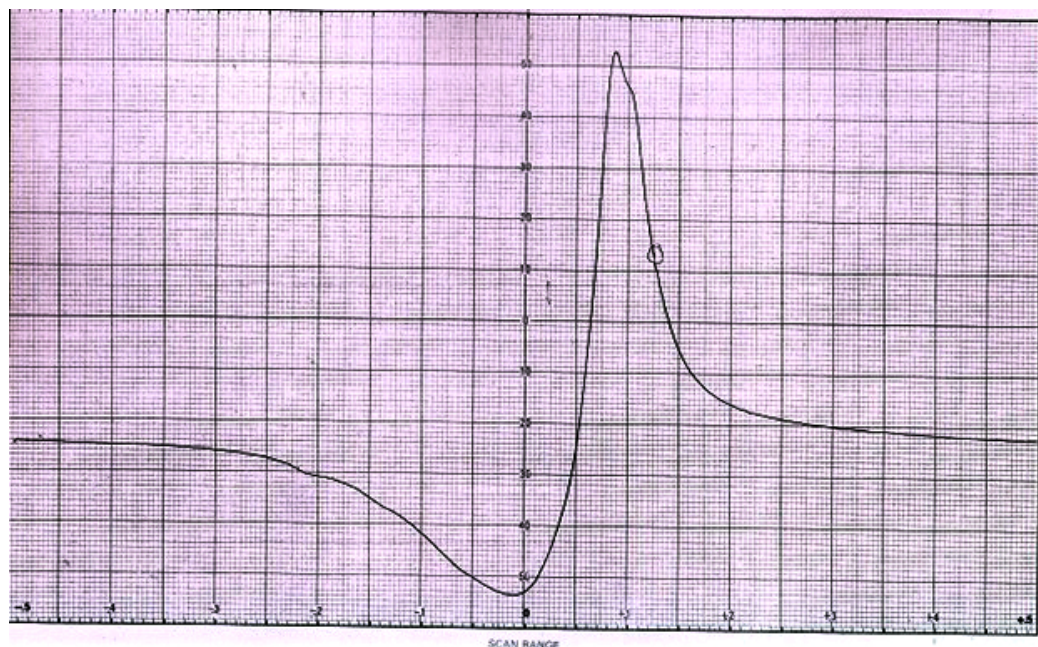
Fig. (5). ESR spectra of complex CC_{pl}AMB.



Fig. (6). ESR spectra of complex $CC_{pr}AMB$.



Fig. (7). ESR spectra of complex $CL_{Cu}AMB$.

4.4. Fungicidal Activities

The antifungal activities of the pure ligands, pure soaps and their corresponding complexes have been evaluated by testing against *Alternaria-alternata* [25 - 26] at a different concentration by Agar plate technique. A perusal of Fig. (8) reveals that all complexes show higher activity than pure soaps and ligands suggesting that complexes are more powerful antifungal agents and benzothiazole and other N, S, O etc [27, 28]. containing compounds are able to enhance the performance of copper soaps. The organic compounds containing amino group play an important role in biology, as it constitutes the repeating unit of the polypeptide macromolecules. These results show that the Cu (II) soap-complexes of ligands are much more toxic than the ligands themselves [29]. The enhanced activity of newly synthesized complexes as compared to those of the ligand can possibly be explained on the basis of chelate formation, the presence of donor atoms, basicity as well as the structural compatibility with molecular nature of the toxic moiety [30]. The enhanced biological activity of complexes is in accordance with the chelation theory. The results of ANOVA [31] for

the antifungal activities for all soaps complexes are shown in Table 5. The predicted R^2 was observed to be in reasonable agreement and closer to 1.0. This confirms that the experimental data are satisfactory. The descriptive statics results shown in Table 6 confirm the satisfactory results in triplet [32].

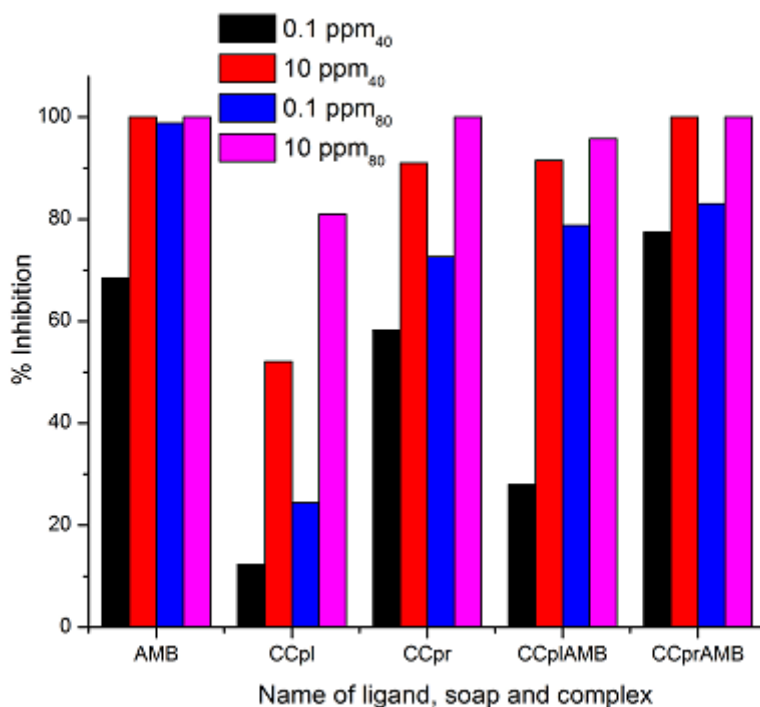


Fig. (8). Antifungal activity of ligand soap and complex (methyl) for fungi *Alternaria Alternata*.

Table 5. Descriptive statics results for antifungal activities of Cu (II) soaps and its complexes.

Compound	conc. (ppm)	Count	Average % Inhibition	Variance	Coefficient Variance	Std. Deviation	Std. Error
CCplS ₄₀	0.1	3	12.4	0.015	0.12	0.072	1.007
	10	3	52.22	0.012	0.11	0.063	0.21
CCplS ₈₀	0.1	3	24.36	0.043	0.2	0.12	0.854
	10	3	80.5	0.16	0.04	0.23	0.496
CCprS ₄₀	0.1	3	58.2	0.01	0.17	0.1	0.057
	10	3	90.76	0.023	0.16	0.152	0.088
CCprS ₈₀	0.1	3	72.55	0.013	0.15	0.115	0.066
	10	3	99.99	0.015	0.01	0.01	0.005
AMB ₄₀	0.1	3	68.23	0.013	0.16	0.115	0.066
	10	3	99.99	0.01	0.1	0.101	0.058
AMB ₈₀	0.1	3	98.8	0.002	0.05	0.051	0.028
	10	3	99.99	1.00E-04	0.01	0.011	0.005
CCplAMB ₄₀	0.1	3	27.99	0.01	0.36	0.102	0.058
	10	3	91.43	0.033	0.19	0.182	0.105
CCplAMB ₈₀	0.1	3	78.74	0.029	0.21	0.172	0.099
	10	3	95.67	0.012	0.11	0.111	0.064
CCplAMB ₄₀	0.1	3	77.4	0.01	0.13	0.1	0.058
	10	3	99.94	0.01	0.1	0.103	0.06
CCplAMB ₈₀	0.1	3	83.27	0.002	0.05	0.046	0.026
	10	3	99.99	3.33E-05	0.005	0.005	0.003

Table 6. ANOVA results for antifungal activities of Cu (II) soaps and its complexes.

Name of compound	SS	df	MS	F	P-value	F crit.	R- SQUARE
CCplS	8408	3	2802	445508	3.16E-21	4.06	0.992
CCprS	3144	3	1048	96521	1.43E-18	4.06	0.995
AMB	2212	3	737	112762	7.7E-19	4.06	0.999
CCplAMB	8734	3	2911	135740	3.67E-19	4.06	0.994
CCprAMB	1207	3	402	69685	5.28E-18	4.06	0.998

CONCLUSION

In conclusion, it can be said that the appearance of enhanced activity may be due to synergistic mechanism *i.e.* the free ligands are less active but on complexation, show more activity in combination with copper (II) soaps. Thus it is evident that concentration plays a vital role in increasing the degree of inhibition. So fungicidal screening data revealed that at lower concentration, the inhibition of growth is less as compared to higher concentration. In all complexes, toxicity increases with the domination of methanol in the solvent mixture of methanol and benzene [40% methanol–benzene and 80% methanol–benzene]. Suggesting that complexes are more effectively playing their role in determining the growth of fungi in the domination of polar solvent in the ternary system: complex + methanol + benzene. Some recent studies about antimicrobial activities of newly synthesized 1, 4-benzothiazines possessing thiazolyl / imidazolyl moieties also support our findings.

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ETHICS APPROVAL AND CONSENT TO PARTICIPATE

Not applicable.

HUMAN AND ANIMAL RIGHTS

No Animals/Humans were used for studies that are base of this research.

CONSENT FOR PUBLICATION

Not applicable.

CONFLICT OF INTEREST

The authors confirm that this article content has no conflict of interest.

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Declared none.

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