

SYLLABUS

MANONMANIAM SUNDARANAR UNIVERSITY, TIRUNELVELI-12

PG - COURSES – AFFILIATED COLLEGES

Course Structure for

M.Sc. Chemistry

(Choice Based Credit System)

(with effect from the academic year 2017- 2018 onwards)

Semester-III				
Part	Subject Status	Subject Title	Subject Code	Credit
	Core - 13	Organic Chemistry – III	PCHM31	4
	Core - 14	Inorganic Chemistry – III	PCHM32	4
	Core - 15	Physical Chemistry – III	PCHM33	4
	Core - 16	Scientific - Research Methodology	PCHM34	4
	Core - 17 Practical - 7	Organic Chemistry Practical – III	PCHL31	2
	Core - 18 Practical - 8	Inorganic Chemistry Practical – III	PCHL32	2
	Core - 19 Practical - 9	Physical Chemistry Practical – III	PCHL33	2

ORGANIC CHEMISTRY – III

Objectives:

- To understand the Aliphatic Nucleophilic substitutions, concept of NMR, Mass Spectroscopy.
- To understand the photochemistry, pericyclic and Hetero cyclic reactions.

Unit-I

Aliphatic nucleophilic substitution and Elimination Reactions:

Aliphatic nucleophilic substitution : Mechanism of S_N1 , S_N2 , S_Ni , S_N1' , S_N2' and S_Ni' reactions- Effect of substrate, nucleophile, leaving group and solvent on the rate of substitution- Ambident nucleophile- NGP- Mechanism of esterifications and ester hydrolysis ($B_{AC}2$ and $A_{AC}2$ mechanisms only)

Elimination reaction: E_1 , E_2 and E_1CB mechanisms- Factors influencing elimination reactions- Hofmann and Saytzeff rules- Pyrolytic elimination- Chugaev and cope reactions-competition between substitution and elimination reactions.



Unit – II: NMR SPECTROSCOPY

^1H -NMR spectroscopy: Basic Principle – number of signals – chemical shift – Factors influencing chemical shift - spin–spin splitting–Proton exchange reactions - classification of spin systems – analysis of AX, AMX and ABX systems – Geminal, Vicinal and long range couplings–NOE in stereochemistry – FTNMR.

C-13 spectroscopy: Principle of proton decoupled and OFF- resonance decoupled C-13 spectroscopy - comparison with ^1H NMR - chemical shifts (aliphatic, olefinic, alkylic, aromatic and carbonyl compounds)

2D NMR spectroscopy: $\text{H}^1\text{--H}^1\text{COSY}$, $\text{H}^1\text{--C}^{13}\text{ COSY}$, NOESY, DEPT and INADEQUATE spectra.

Unit – III:**MASS SPECTROSCOPY**

Basic Principles– Base peak – molecular ion – nitrogen rule – metastable ions – isotopic peak - daughter ions – Mc–Lafferty rearrangement – RDA – General rules for fragmentation pattern – Fragmentation pattern of simple compounds of hydrocarbons, alcohols, amines, aldehyde, ketone, ether, acids, phenols ,nitro compounds, alicyclic compounds .

Alternative electron impact ionization technique– CI, FAB, ESI – MS, MALDI –MS, MALDI-TOF , ICP- MS. One conjunction problem based on UV, IR, ^1H NMR, ^{13}C NMR and Mass spectroscopic techniques is compulsory under section – c. Problems shall be based on the reference books.

Unit-IV :**Organic photochemistry and pericyclic reactions**

Organic photochemistry: Jablonskii diagrams-intersystem crossing-energy transfer process- Photosensitization- alpha cleavages or Norrish type-I and Norrish Type II cleavages - Paterno-Buchi reactions- Barton reaction, cis-trans isomerisation. - Di- π methane rearrangement.

pericyclic reactions:

Atomic and molecular orbitals-Woodward-Hoffmann rules, FMO and correlation diagram approaches:

Electrocyclic reaction: con and dis rotatory motions for $4n$ and $4n+2$ system (butadiene and 1,3,5-hexatrienes)- Stereochemical course of electro cyclic reaction in terms of conservation of orbital symmetry.

Cycloaddition: suprafacial and antarafacial, $[2+2]$ and $[4+2]$ cyclo addition reactions (ethylene and butadiene)

Sigmatropic rearrangements - $[i,j]$ shift of C-H and C-C bonds (1,3) and (1,5) carbon migration.

Unit-V :**Heterocyclic and biomolecules**

Synthesis and reactions of indole, oxazole, imidazole, thiazole, Reserpine and quinine chromans, pyrimidine, pyridazine, pyrazine, coumarins, benzopyrones and anthocyanins-synthesis of flavones, isoflavones, flavonol, and quercetin -Biosynthesis of flavonoids. Synthesis



Pyranose and furanose forms of aldohexose and ketohexose-methods used for the determination of ring size-A Detailed study on the structure of maltose, lactose and starch.

REFERENCES

1. J. March, 'Advanced organic chemistry', Fourth Edition, John Wiley and Sons, New York, 2006.
2. Depuy, E.C.H. and Chapman, O.S., "Molecular reactions and photochemistry", Prentice Hall, New York, 1988.
3. I.L. Finar, 'Organic Chemistry', volume 2, sixth Edition, Pearson Education Inc., Singapore, 2006.
4. Raj K. Bansal, 'Organic Reaction mechanisms', Tata Mc Graw Hill, Third Edition, 2007
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9. F.A. Carey and J. Sundberg, 'Advanced Organic chemistry' part A and B, Plenum Press, 2005.
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15. E.S. Gould, 'Mechanism and structure in organic chemistry' Holt, Rinehart and Winston Inc., 1959
16. F.A. Carey, Organic chemistry – Tata Mc Graw Hill, Delhi, 5th edition 2005.
17. Stryer, L., "Biochemistry", Fifth edition, W.H. Freeman and company, San Francisco, 2002.
18. Jain, J.L., "Fundamentals of Biochemistry", Fourth edition, S. Chand & Company Limited, New Delhi. 2007
19. Bansal, K., "Heterocyclic Chemistry", Fourth edition, New Age International, New Delhi, 2005.



INORGANIC CHEMISTRY - III

Objectives:

- To introduce organometallic compounds and to study their catalytic applications in homogeneous and heterogeneous systems.
- To study the applications of NMR and EPR techniques in inorganic systems.
- To understand the basic principles and applications of thermo and spectro analytical techniques.
- To introduce inorganic photochemistry and to study applications in various systems.

Unit I

ORGANOMETALLIC CHEMISTRY – I

The 18 e⁻ and 16 e⁻ rules and its correlation to stability – Synthesis and structures of metal carbonyls, metal nitrosyls and dinitrogen complexes – Substitution reactions of metal carbonyls - IR spectral applications – identifications of bridging and terminal CO groups – Stretching mode analysis of metal carbonyls – evidence for M-M bonds. Synthesis, properties and structural features of metal complexes with alkene, alkyne, allyl and arene systems. Metallocenes – synthesis, properties, structure and bonding with particular reference to ferrocene and zirconocene – covalent versus ionic bonding in zirconocene. Template synthesis of macrocyclic ligands.

Unit II

ORGANOMETALLIC CHEMISTRY – II

Organometallic compounds as catalysts and the requirements: Agostic interaction – Oxidative addition and reductive elimination - insertion and elimination reactions – nucleophilic and electrophilic attack of coordinating ligands - cyclometallation reactions. Homogeneous catalysis: Wilkinson's catalyst and hydrogenation reactions, Tolman catalytic loop; hydroformylation (oxo) reaction, Wacker and Monsanto acetic acid processes. Cluster compounds, polymer-supported and phase-transfer catalysis. Heterogeneous catalysis: synthesis gas and water gas shift reactions; Fischer Tropsch process and synthetic gasoline, Ziegler-Natta polymerization and mechanism of stereoregular polymer synthesis. Cyclooligomerisation of acetylenes (Rupe's or Wilke's catalyst) – Olefin isomerisation using Ni catalyst – olefin metathesis catalysed by Schrock type carbene.

UNIT-III:

SPECTRAL METHODS TO THE STUDY OF INORGANIC COMPOUNDS – I

NMR SPECTROSCOPY: ³¹P, ¹⁹F and ¹⁵N – NMR – applications in structural problems based on number of signals, multiplicity, anisotropy (like H₃PO₃, H₃PO₂, [HNi(PPh₃)₄]⁺, SF₄, TiF₄, PF₅, HPF₂, H₂PF₃, PF₃(NH₂)₂, P₄S₃, P₄N₄Cl₆(NHC₆H₅)₂, P₃N₃(CH₃)₂C₁₄, NF₃, NH₃ – mer- and fac-Rh(PPh₃)₃C₃, fluxional molecules (including organometallic compounds) and study of fluxionality by NMR technique - NMR of paramagnetic molecules - contact shifts. Evaluation of rate constants -



monitoring the course of reaction using NMR.

EPR spectroscopy: Factors affecting magnitude of g-values - Zero field splitting and Kramers' degeneracy - Application of EPR in the study of transition metal complexes based on number of signals, multiplicity, anisotropy (bis(salicylaldimine)copper(II), $[\text{Cu}(\text{bpy})_3]^{2+}$, $[\text{Cu}(\text{Phen})\text{Cl}_2]$, $[(\text{NH}_3)_5\text{Co}-\text{O}_2-\text{Co}(\text{NH}_3)_5]^{5+}$, $\text{Co}_3(\text{CO})_9\text{Se}$, $\text{Co}_3(\text{CO})_9\text{Rh}$, $[\text{CoF}_6]^{4-}$, $[\text{CrF}_6]^{3-}$, $\text{VO}(\text{acac})_2$, $[\text{VO}(\text{H}_2\text{O})_6]^{2+}$, $[\text{Fe}(\text{CN})_5\text{NO}]^{2-}$). Applications in predicting the covalent character of M-L bond and Jahn-Teller distortion in Cu(II) complexes. EPR spectroscopy of metallobiomolecules: copper and iron proteins.

UNIT – IV:

THERMOANALYTICAL AND SPECTROANALYTICAL METHODS

Theory and principles of thermogravimetric analysis, differential thermal analysis and differential scanning calorimetry – characteristic features of TGA and DTA curves – factors affecting TGA and DTA curves – complementary nature of TGA and DTA – applications of thermal methods in analytical chemistry – thermometric titrations – the study of minerals and metal compounds. Principle and applications of spectrophotometry, spectrofluorimetry, atomic absorption spectroscopy and atomic emission spectroscopy based on plasma sources.

UNIT -V :

PHOTOCHEMISTRY OF METAL COMPLEXES

Frank Condon and thermally equilibrated excited (THEXI) states – properties of excited states of metal complexes – types of excited states, photophysical processes: bimolecular deactivation and energy transfer, photochemical processes: electron transfer reactions, isomerisation and substitutional processes – Photochemistry of Cr(III) and Co(III) complexes – Photophysical and photochemical properties of $[\text{Ru}(\text{bpy})_3]^{2+}$. Applications of inorganic photochemistry: photochemical conversion and storage of solar energy – inorganic photochemistry at semi-conductor electrodes - Catalyzed photoreduction of CO_2 and CO – TiO_2 as a green photocatalyst in removing air and water pollutants.

REFERENCES

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2. F. Albert Cotton, Geoffrey Wilkinson, Carlos A. Manic and Manfred Bochman, Advanced Inorganic Chemistry, Wiley Interscience Publication, 6th Edition, 1999.
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4. R. H. Crabtree, The Organometallic Chemistry of the Transition Metals, 4th Edition, John Wiley, 2005.
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 12. David W. H. Rankin, Norbert W. Mitzel, Carole A. Morrison, Structural Methods in Molecular Inorganic Chemistry, John Wiley & Sons, Ltd, 1st Edition, 2013.
 13. Robert A. Scott and Charles M. Lukehart, Applications of Physical Methods to Inorganic and Bioinorganic chemistry, John Wiley & Sons Ltd, 2007.
 14. Philip H Rieger, Electron Spin Resonance Analysis and Interpretation, The Royal Society of Chemistry, 2007.
 15. John A. Weil, James R. Bolton, Electron Paramagnetic Resonance Elementary Theory and Practical Applications, 2nd Edition, John Wiley & Sons, 2007.
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 24. V. Balzani and A. Juris, Coord. Chem. Rev., 211, 97–115, 2001.



PHYSICAL CHEMISTRY- III

Objective:

- Learning the concepts of Group Theory
- To understand the Principles and applications of various spectroscopy

UNIT-I:

Group Theory-I

Symmetry elements and operations. Group Postulates and types of groups. Identification of Point groups of molecules and Schoenflies symbols. Construction of multiplication table for C_{2v} , C_{3v} and C_{2h} . Sub-groups and classes of symmetry operations. Rule of similarity transformations. Matrix representations of symmetry operations. Use of atomic wave functions as bases for point group representations. Reducible and irreducible representations. The Great Orthogonality theorem. Properties of Reducible and irreducible representations. Construction of character tables for C_{2v} , C_{3v} , C_{4v} , C_{2h} , and D_2 point groups by using The Great Orthogonality theorem.

UNIT-II :

Group Theory -II :

Standard Reduction Formula, Vibrational modes as bases for group representations-Normal mode analysis for non linear molecules H_2O , $POCl_3$, trans- N_2F_2 and $PtCl_4$. Symmetry selection rules for infrared and Raman spectra. Mutual exclusion principle. Determination of Hybridisation of atomic orbitals in non-linear molecules (CH_4 , XeF_4 , and PF_5). Electronic spectra of ethylene and formaldehyde molecules. Construction of Projection operators and Molecular orbitals by Symmetry Adapted Linear Combinations. Simplification of HMO calculations using group theory. Calculation of delocalization energy for ethylene, trans-1,3 –butadiene, and benzene systems.

UNIT – III:

Nuclear Magnetic Resonance Spectroscopy

Theory of Proton NMR spectroscopy, Chemical shift and its measurement, Factors influencing chemical shift, Solvents used in NMR, solvents shift-concentration and temperature effects-hydrogen bonding. Theory of Spin-spin splitting-Magnitude of coupling-coupling constants, J, First-order spectra of complex systems, chemical and magnetic equivalence in NMR, Proton exchange reactions, Factors influencing coupling constant, J. Theory and Principle of ^{13}C , ^{19}F , ^{31}P NMR-Range of chemical shift values, spectra of typical examples. FT NMR-FIDs. Theory of Spin-spin splitting and double irradiation, InterNuclear Double Resonance (INDOR) and Selective Population Inversion (SPI), Nuclear OverhauserEffect (NOE), 2D NMR-shift correlation spectra-COSY, Magnetic Resonance Imaging (MRI).

UNIT-IV:

NQR and EPR spectroscopy



Electron paramagnetic resonance spectroscopy: theory of EPR spectroscopy, presentation of the spectrum, nuclear hyperfine splitting in isotropic systems. EPR spectra of anisotropic systems: anisotropy in g-value, causes of an isotropy, anisotropy in hyperfine coupling. Double resonance in ESR, Zero field splitting and Kramers' degeneracy.

Theory and Principle of NQR spectroscopy-Nature of electric field gradient, Energy levels and selection rules, Interaction of electric quadrupole with electromagnetic radiation, nuclear orientations, the asymmetry parameter, quadrupole transitions in spherical, axially symmetric fields and not axially symmetric fields. Applications of NQR spectra.

UNIT-V:

Electronic Spectroscopy, Mossbauer Spectroscopy and Mass Spectrometry

Electronic Spectroscopy-Electronic Spectrum of diatomic molecules-Born-Oppenheimer approximation, Progressions, Franck-Condon Principle, Dissociation Energy and dissociation products, Rotational Fine structure of Electronic-Vibration Transitions, The Fortrat diagram, Predissociation, Electronic states of atoms, Electron orbitals in diatomic molecules, Electronic states of diatomic molecules, Potential energy curves for Electronic states of diatomic molecules.

Photoelectron Spectroscopy-Basic Principles, Ultra-Violet Photoelectron Spectroscopy, X-ray Photoelectron Spectroscopy, Chemical information from Photoelectron Spectroscopy. Mössbauer spectra: Theory and Principle of Mössbauer spectra, isomer shift, quadrupole interactions, magnetic hyperfine interaction, Doppler shift, recoil energy, experimental technique-sources, absorber, calibration, Chemical applications.

Mass spectrometry: Operation and representation of spectra. Effect of combination of high energy electron with a molecule. Finger print application and the interaction of mass spectra, Effect of isotopes on the appearance of a mass spectrum, Molecular weight determinations.

REFERENCE BOOKS

1. Introductory Group Theory For Chemists- George Davidson
2. F. Albert Cotton, Chemical Applications of Group Theory, Third Edition John Wiley & Sons, Singapore 2003.
3. V. Ramakrishnan and M. S. Gopinathan: Group Theory in chemistry, Vishal Publication, 1986.
4. Robert L. Carter, Molecular Symmetry and Group Theory John Wiley and Sons, Inc., New York, 1998.
5. R.L. Flurry, Jr, Symmetry Groups – Prentice Hall, New Jersey 1980.
6. Group Theory and its applications to Chemistry by K. V. Raman
7. Group Theory and Spectroscopy by K. Veera Reddy
8. Group Theory and its Chemical applications by B.K. Battacharya
9. C.N. Banwell and E. M. McCash, Fundamentals of Molecular Spectroscopy, 4th ed., Tata McGraw Hill, New Delhi, 2000.
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13. Spectroscopy (Atomic and Molecular) Gurdeep R. Chatwal and Sham K. Anand, Himalaya Publishing House
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16. Introduction to Molecular spectroscopy, G.M. Barrow, McGraw-Hill international editions.
17. R.M. Silverstein and F. X. Webster, Spectroscopic Identification of Organic Compounds, 6th ed., John Wiley & Sons, New York, 2003.



SCIENTIFIC - RESEARCH METHODOLOGY

Objectives:

- How to learn the survey for literature, chemical abstract, choosing a research problem and scientific writing, characterization and data analysis, computer searches and literature.
- How to apply for the various finding agencies.

Unit – I:

LITERATURE SURVEY

Source of chemical information – primary, secondary, tertiary sources- literature survey-Indexes and abstracts in science and technology – Applied science and technology index, chemical abstracts, chemical titles, current chemical reactions, current contents and science citation index.

Classical and comprehensive reference works in chemistry-synthetic methods and techniques, treatises, reviews, patents and monographs.

UNIT - II :

CHEMICAL ABSTRACTS:

Current awareness searching: CA weekly issues, CA issue indexes. Retrospective searching: CA volume indexes-general subject index, chemical substance index-formula index, index of ring systems, author index, patent index. CA collective indexes: collective index (CI), decennial index (DI). Access points for searching CA indexes- Index guide, general subject, terms, chemical substance names, molecular formulas, ring systems, author names, patent numbers. Locating the reference: finding the abstract, finding the original document chemical abstract - service source index.

UNIT –III:

CHOOSING A RESEARCH PROBLEM AND SCIENTIFIC WRITING

Identification of research problem – assessing the status of the problem - guidance from the supervisor – actual investigation and analysis of experimental results – conclusions.

Scientific writing-research reports, thesis, journal articles and books.

Steps to publishing a scientific article in a journal – types of publications-communications, articles, reviews, when to publish, where to publish ,specific format required for submission.

Documenting- Abstracts-indicative (or) descriptive abstracts, informative abstract, footnotes, end notes, referencing styles-bibliography-journal abbreviations (CASSI), abbreviation used in scientific writing.

Unit –IV:

INSTRUMENTAL CHARACTERIZATION AND DATA ANALYSIS

Principle and Sample preparation of UV, FT-IR, TEM, SEM, EDAX, AFM and XRD characterization of observed results – Data analysis - Report.



Errors in chemical analysis – classification of errors – determination of accuracy of methods – improving accuracy of analysis – significant figures – mean, standard deviation – comparison of results : “t” test, “f” test, Q test and “chi” square test – rejection of results – presentation of data.

UNIT –V:

COMPUTER SEARCHES AND LITERATURE

ASAP – Alerts, CA Alerts, scifinder, chemport, science direct, STN international, journal home pages. Online browsing of research articles – online submission of research papers in various Journals (ACS, RSC, Elsevier, Springer etc.) –Instructions to the authors – Impact factors. Writing project proposal to funding agencies (UGC, DST etc.).

REFERENCES

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2. A.J. Durston, Thesis and assignment writing.
3. R.O.Bullet, Preparing thesis and other manuscripts.
4. R. L. Dominoski, Research Methods, Prentice Hall, 1981.
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9. J. S. Dodd, Ed., The ACS Style Guide: A Manual for Authors and Editors; American Chemical Society: Washington, DC, 1985.
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13. www.ugc.ac.in/pdfnews/7716504_12th-plan-guidelines.pdf
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18. B.K. Sharma, Instrumental methods of chemical analysis.
19. D.A. Skoog and M. West, Fundamentals of analytical chemistry.
20. J.D. Dick, Analytical chemistry.
21. S.M. Khopkar, Basic concepts of analytical chemistry.



Organic Chemistry Practical – III

Estimations, two stage preparations and Spectral interpretation have been included as the practical components.

Microscale preparations are recommended for the simple reason, they are both economic-friendly and eco-friendly

A. List of Estimations

1. Ethylmethylketone
2. Acetone
3. Saponification value of an oil
4. Determination of Percentage purity in an unsaturated acid.
5. Estimation of hydroxyl group

B. List of Two stage preparations

1. Benzaldehyde Benzoic acid m-nitro benzoic acid
2. Acetanilide p-acetanilide p-Bromoaniline
3. Methyl benzoate m-nitro methyl benzoate m-nitro benzoic acid
4. Acetanilide p-nitro acetanilide p - nitroaniline
5. Benzophenone Benzo phenone oxime Benzanilide

Students are expected to submit the recrystallised samples of the final products, at the time of practical examination, for evaluation by the examiners.

C. For Class work Only :

1. Download the following spectra from internet and give interpretation.

Differentiate the following pair by ¹H NMR spectra

- (a) Maleic acid and Fumaric acid.
- (b) Aqueous ethyl alcohol and Pure ethyl alcohol.
- (c) Dimethyl Ether and Aqueous ethyl alcohol.

Interpret the following C-13 NMR Spectra.

- (a) OFF- Resonance decoupled C-13 spectrum of menthol.
- (b) DEPT spectrum of isopentyl acetate.
- (c) INADEQUATE spectrum of 2- butanone.

Interpret the mass spectrum of anisole and benzoic acid.

N.B: 1. Section -C is course work only.

2. It is for the purpose of internal assessment only.

REFERENCES

1. F.C.Mann and B.C.Saunders, Practical organic chemistry, Fourth edition, ELBS, 1970
2. A.I. Vogel, A Text book of Practical organic chemistry.
3. A.I. Vogel, A Text book of Quantitative Organic Analysis, 1989.
4. Raj K. Bansal, Laboratory Manual of Organic Chemistry, Second Edition, Wiley Eastern Ltd., 1990



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8. V.K.Srivastava and K.K.Srivastava, Introduction to Chromatography-Theory and Practice, S.Chand & Co., 1987.



INORGANIC CHEMISTRY PRACTICAL – III

I. Quantitative estimation of a mixture containing two metal ions (Volumetric and Gravimetric Estimations).

1. Estimation of mixture of Cu^{2+} and Ni^{2+} ions.
2. Estimation of mixture of Cu^{2+} and Zn^{2+} ions.
3. Estimation of mixture of Fe^{2+} and Cu^{2+} ions.
4. Estimation of mixture of Fe^{2+} and Ni^{2+} ions.
5. Estimation of mixture of Ca^{2+} and Mg^{2+} ions.
6. Estimation of mixture of Ca^{2+} and Ba^{2+} ions.

II. Analysis of ores and alloys (course work).

III. One day visit to Industry/Research Institution and submission of a minor report.

REFERENCES

1. G.H. Jeffery, J. Bassett, J. Mendham and R.C. Denney, Vogel's Textbook of Quantitative Chemical Analysis, Revised 5th edition, ELBS, 1989.
2. Mounir A. Malati, Experimental Inorganic/Physical Chemistry - An Investigative, Integrated Approach to Practical Project Work, Woodhead Publishing Limited, Reprint 2010



PHYSICAL CHEMISTRY PRACTICAL-III

Objective:

- To learn and apply the Principles of Potentiometric Titrations.
- To understand the Principles and applications of Adsorption

I. POTENTIOMETRIC TITRATIONS

- (a) Acid alkali titrations.
- (b) Redox titrations (i) Fe^{2+} vs $\text{Cr}_2\text{O}_7^{2-}$
- (d) Determination of dissociation constant of weak acids.
- (e) Determination of activity and activity coefficient of ions.
- (f) Determination of pH of a buffer solution using a quinhydrone electrode.

II. TITRATION USING PH METER

- (a) Determination of dissociation constant of Weak acid.

III. ADSORPTION

Freundlich Adsorption isotherm

Adsorption of oxalic acid on charcoal.

REFERENCES

1. J.B.Yadav, "Advanced Practical Physical chemistry", 20th Edn., GOEL publishing House, Krishna Pakashan Media Ltd., (2001).
2. Findlay's "Practical Physical Chemistry" Revised and edited by B.P. Levitt 9th Edn., Longman, London, 1985.
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